SITE INVESTIGATION REPORT

VOLUME II

FINAL

ALABAMA AIR NATIONAL GUARD 187th FIGHTER GROUP

DANNELLY FIELD MUNICIPAL AIRPORT MONTGOMERY, ALABAMA

NOVEMBER 1995



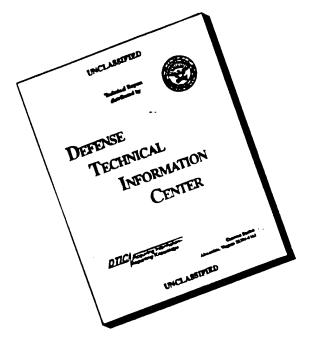
Hazardous Waste Remedial Actions Program
Environmental Restoration and Waste Management Programs

Oak Ridge, Tennessee 37831-7606

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SOIL GAS/GROUNDWATER SURVEY RESULTS

SOIL GAS AND GROUND WATER SURVEY ALABAMA AIR NATIONAL GUARD MONTGOMERY, ALABAMA



RGET ENVIRONMENTAL SERVICES, INC.

SOIL GAS AND GROUND WATER SURVEY ALABAMA AIR NATIONAL GUARD DANNELLY FIELD POL FACILITY MONTGOMERY, ALABAMA

PREPARED FOR

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MARCH 1991

EXECUTIVE SUMMARY

On February 19-21, 1991, TARGET Environmental Services, Inc. (TARGET) conducted a soil gas and ground water survey at Alabama Air National Guard Dannelly Field Petroleum, Oil and Lubricant (POL) Facility in Montgomery, Alabama, where underground storage tanks containing JP-4 may have leaked and minor fuel spillage has been reported. Soil gas and ground water headspace samples were analyzed by GC/FID for petroleum hydrocarbons and by GC/ECD for the chlorinated hydrocarbons, trichloroethene (TCE) and tetrachloroethene (PCE).

Elevated levels of Total Volatiles were present in all soil gas and ground water samples. Toluene, ethylbenzene and the xylene isomers were present in all ground water and most soil gas samples. Benzene was the most restricted of the standardized petroleum analytes, particularly in the soil gas phase. The chlorinated hydrocarbons, PCE and TCE, were not present above their detection limit in any of the soil gas or ground water samples collected from the POL facility.

Introduction

CH2M HILL contracted Target Environmental Services, Inc. (TARGET) to conduct a soil gas and ground water survey at the Petroleum, Oil and Lubricant (POL) Facility, Alabama Air National Guard Dannelly Field, Montgomery, Alabama, as part of a site investigation. There is evidence of past leakage from JP-4 underground storage tanks and minor JP-4 spillage has been reported. The purpose of the survey is to aid in the placement of soil borings and monitoring wells. The field phase of the soil gas survey was conducted on February 19-21, 1991.

Detectability

The soil gas survey data presented in this report are the result of precise sampling and measurement of contaminant concentrations in the vadose zone. Analyte detection at a particular location is representative of vapor, dissolved, and/or liquid phase contamination at that location. The presence of detectable levels of target analytes in the vadose zone is dependent upon several factors, including the presence of vapor-phase hydrocarbons or dissolved or liquid concentrations adequate to facilitate volatilization into the unsaturated zone.

Terminology

In order to prevent misunderstanding of certain terms used in this report, the following clarifications are offered:

The term "feature" is used in reference to a discernible pattern in the contoured data. It denotes a contour form rather than a definite or separate chemical occurrence.

The term "occurrence" is used to indicate an area where chemical compounds are present in sufficient concentrations to be detected by the analysis of soil vapors. The term is not indicative of any specific mode of occurrence (vapor, dissolved, etc.), and does not necessarily indicate or suggest the presence of "free product" or "phase-separated hydrocarbons."

The term "anomaly" refers to an area where hydrocarbons were measured in excess of what would normally be considered "natural" or "background" levels.

The term "analyte" refers to any of the hydrocarbons standardized for quantification in the chromatographic analysis.

The term "vadose zone" represents the unsaturated zone between the ground water table and the ground surface.

The term "indicates" is used when evidence dictates a unique conclusion. The term "suggests" is used when several explanations of certain evidence are possible, but one in particular seems more likely. As a result, "indicates" carries a higher degree of confidence in a conclusion than does "suggests."

Field Procedures

Ground water samples were collected at a total of 10 locations at the site, as shown in Figure 1A. To collect the samples a slide hammer was used to advance connected 3' sections of 1" diameter threaded steel casing down to a depth of 6'. Sample 27-W was collected at a depth of 5'. Water was allowed to fill the pipe. A stainless steel bailer was used to collect the water samples. Samples were placed in amber glass vials, sealed, labeled and taken to the mobile laboratory.

A field control sample of water was collected at the beginning of each day's field activities and after the fourth and eighth samples on the second day. This QA/QC sample was obtained using distilled water.

Soil gas samples were collected at a total of 12 locations at the site, as shown in Figure 1B. To collect the samples, a vanmounted hydraulic probe was used to advance connected 3' sections of 1" diameter threaded steel casing down to a depth of 15'. The entire sampling system was purged with ambient air drawn through an organic vapor filter cartridge. A teflon line was inserted into the casing to the bottom of the hole, and the bottom-hole line perforations were isolated from the up-hole annulus by an inflatable packer. A sample of in-situ soil gas was then withdrawn through the probe and used to purge atmospheric air from the sampling system. A second sample of soil gas was withdrawn through the probe and encapsulated in a pre-evacuated glass vial at two atmospheres of pressure (15 psig). The self-sealing vial was detached from the sampling system, packaged, labeled, and taken to

the mobile laboratory for analysis.

Field control samples of soil gas were collected at the beginning and end of each day's field activities. These QA/QC samples were obtained by drawing ambient air through a dust and organic vapor filter cartridge and collecting in the same manner as described above.

Prior to the day's field activities all sampling equipment, slide hammer rods, steel casing and bailers were decontaminated by washing with soapy water and rinsing thoroughly. Internal surfaces were flushed dry using pre-purified nitrogen or filtered ambient air, and external surfaces were wiped clean using clean paper towels.

Laboratory Procedures

All analyses were performed on site in TARGET's climate controlled mobile laboratory.

Headspaces of ground water and aqueous standards were prepared by pipetting equal volumes of liquid from the sample containers. Sample containers were placed in a heating block at 75-80°C for 5 minutes prior to injection of a vapor sample from the headspace of the sample container.

All of the soil gas and ground water headspace samples collected during the field phase of the survey were subjected to dual analyses. One analysis was conducted according to EPA Method 601 (modified) on a gas chromatograph equipped with an electron capture detector (ECD), but using direct injection instead of purge and trap. Specific analytes standardized for this analysis were:

trichloroethene (TCE) tetrachloroethene (PCE)

The chlorinated hydrocarbons were chosen because of their common usage in industrial solvents.

The second analysis was conducted according to EPA Method 602 (modified) on a gas chromatograph equipped with a flame ionization detector (FID), but using direct injection instead of purge and trap. The analytes selected for standardization in this analysis were:

benzene
toluene
ethylbenzene
meta- and para- xylene
ortho-xylene

These compounds were chosen because of their utility in evaluating the presence of fuel products, or petroleum based solvents.

The FID Total Volatiles values were generated by summing the areas of all chromatogram peaks and calculated using the instrument response factor for toluene. Injection peaks, which also contain the light hydrocarbon methane, were excluded to avoid the skewing of the Total Volatiles (Totals) values due to injection disturbances and biogenic methane. For samples with low hydrocarbon concentrations, the calculated Total Volatiles concentration is occasionally lower than the sum of the individual analytes. This is because the response factor used for the Total Volatiles calculation is a constant, whereas the individual analyte response factors vary with concentration. It is important to understand that the Total Volatiles levels reported are relative, not absolute, values.

Separate standard curves were used for soil gas and ground water headspace. The analytical equipment was calibrated using an instrument-response curve and injection of known concentrations of the above standards. Retention times of the standards were used to identify the peaks in the chromatograms of the field samples and their response factors were used to calculate the analyte concentrations. The tabulated results of the laboratory analyses of the soil gas and ground water headspace samples are reported in parts per billion volume to volume (ppb v/v) in Tables 1 and 2.

Map sample points with no data shown indicate that the analyte concentrations in the sample were below the detection limit.

For QA/QC purposes, a duplicate analysis was performed on every tenth field sample. Laboratory blanks of nitrogen gas (99.999%) were also analyzed after the tenth soil gas sample, while

a laboratory blank of distilled water was analyzed after the tenth ground water headspace sample.

Discussion and Interpretation of Results

In order to provide graphic presentation of the results, ground water data sets in Table 1 have been mapped and contoured to produce Figures 2 through 6. Soil gas data sets in Table 2 have been mapped in Figures 7-12. Soil gas data were not contourable.

Analysis of the ground water via headspace indicated elevated levels of Total Volatile hydrocarbons (Figure 2) in all ground water samples. Benzene and toluene concentrations (Figures 3 and 4) were highest in the northern portion of the POL. Benzene was not present in the ground water above the reported detection limit in samples from the southeastern portion of the POL, while toluene concentrations were relatively low in this area. Ethylbenzene (not mapped), m- and p-xylene (Figure 5) and o-xylene (Figure 6) samples collected in concentrations were highest in southeastern portion of the site (Stations 23-W and 26-W). GC/ECD analysis of the ground water headspace samples indicated that tetrachloroethene (PCE) and trichloroethene (TCE) were not present above their respective detection limits in any of the samples.

GC/FID analysis of soil gas samples from the periphery of the POL facility revealed elevated levels of Total Volatiles in all soil gas samples. The highest concentration occurred on the eastern side of the facility north of the building (Station 20). Benzene (Figure 8), the most restricted of the standardized analytes, was present at only two locations on the northeastern side of the facility (Stations 18 and 19). The toluene concentration was highest on the western side of the facility at Station 15, with lower levels in some other samples on the western and

northeastern sides. Ethylbenzene, m- and p-xylene and o-xylene levels (Figures 10-12) were relatively high on the eastern side of the facility at Station 20. A relatively high level of m- and p-xylene also occurred on the western side at Station 14. Elevated concentrations of ethylbenzene and m- and p-xylene were observed at many of the remaining sampling locations, while o-xylene was quite restricted in its occurrence. GC/ECD analysis of the soil gas samples indicated that PCE and TCE were not present above their respective detection limits in any of the soil gas samples.

Elevated levels of Total Volatiles were present in all soil gas and ground water samples. Toluene, ethylbenzene and the xylene isomers were present in all ground water and most soil gas samples. Benzene was the most restricted of the standardized petroleum analytes, particularly in the soil gas samples. The chlorinated hydrocarbons, PCE and TCE, were not present above their detection limits in any of the soil gas or ground water samples collected from the POL facility.

TABLE 1

*ANALYTE CONCENTRATIONS IN GROUND WATER VIA HEADSPACE ANALYSIS CONCENTRATIONS IN PARTS PER BILLION (V/V)

SAMPLE	BENZENE	TOLUENE	ETHYL- BENZENE	m- & p- XYLENE	O- XYLENE	TOTAL VOLATILES	TOE	PCE
23-W	<11	556	20,100	29,290	12,930	361,600	<7.0	0.9>
2 4 – W	<11	384	4,963	4,629	2,963	73,000	<7.0	<6.0
2 1 1 C	153	524	14,460	14,240	4,466	366,900	<7.0	0.9>
M 90	<11	723	25,290	24,920	12,060	522,200	<7.0	0.9>
100 100 100 100 100 100 100 100 100 100	, , , , , , , , , , , , , , , , , , ,	703	6,905	7,028	3,612	231,600	<7.0	0.9>
M=80	, L	3.401	11,530	12,160	6,017	270,400	<7.0	<6.0
M - 00	25.810	10.780	17,230	17,150	7,763	478,000	<7.0	<6.0
M-08	16,670	5,956	10,670	5,374	5,606	308,900	<7.0	0.9>
3 - 1 - 1 - 1 - 1 - 1 - 1 - 1 - 1 - 1 -	13,80		7,048	6,267	2,163	164,400	<7.0	<6.0
32-W	•	105		1,840	1,057	35,450	<7.0	<6.0

^{*} BTEX WERE ANALYZED VIA GC/FID AND HALOCARBONS WERE ANALYZED VIA GC/ECD.

TCE = TRICHLOROETHENE PCE = TETRACHLOROETHENE

¹CALCULATED USING THE SUM OF THE AREAS OF ALL INTEGRATED CHROMATOGRAM PEAKS AND THE INSTRUMENT RESPONSE FACTOR FOR TOLUENE

TABLE 1 (cont)

*ANALYTE CONCENTRATIONS IN GROUND WATER VIA HEADSPACE ANALYSIS CONCENTRATIONS IN PARTS PER BILLION (V/V)

SAMPLE	BENZENE	TOLUENE	ETHYL- BENZENE	m- & p- XYLENE	O- XYLENE	TOTAL VOLATILES1	TCE	PCE
FIELD CON	FIELD CONTROL SAMPLES	ro!						
W- L	<11	<11	<11	<11	<11	<11	<7.0	<6.0
: A	<11	<11	<11	<11	<11	1,636	<7.0	<6.0
- 13 - 13	1 5		16	26	24	2,132	<7.0	<6.0
= 3- - 12- - 13- -	\ 	<11>	<11	15	15	131	<7.0	<6.0
. M	117	<11	<11	<11	<11	141	<7.0	<6.0
LABORATOR	LABORATORY DUPLICATE ANALYSES	ANALYSES						
32 - W	19	105	1.576	1,840	1,057	35,450	<7.0	0.9>
32-WR	5 6 6 8	114	1,974	2,188	1,099	46,630	<7.0	<6.0
LABORATORY BLANKS	Y BLANKS							
BMCHM-1	<11	<11	<11	12	13	407	<7.0	0.9>
				-				

^{*} BIEX WERE ANALYZED VIA GC/FID AND HALOCARBONS WERE ANALYZED VIA GC/ECD.

TCE = TRICHLOROETHENE PCE = TETRACHLOROETHENE

¹CALCULATED USING THE SUM OF THE AREAS OF ALL INTEGRATED CHROMATOGRAM PEAKS AND THE INSTRUMENT RESPONSE FACTOR FOR TOLUENE

TABLE 2

* LABORATORY RESULTS OF SOIL GAS SAMPLES CONCENTRATIONS IN PARTS PER BILLION (V/V)

SAMPLE	BENZENE	TOLUENE	ETHYL- BENZENE	m- & p- XYLENE	o- XYLENE	TOTAL VOLATILES ¹	TCE	PCE
11	<287	<243	<211	<211	<211	2,803	<17	<6.8
12	<287	<243	211	570	<211	33,440	<17	×6.8
13	<287	<243	<211	359	<211	22,320	<17	<6.8
	<287	316	781	2,110	969	75,940	<17	8.9>
	<287	3,645	338	844	359	80,030	<17	<6.8
16	<287	·V	<211	359	<211	7,333	<17	<6.8
	<287	292	253	1,287	232	41,300	<17	<6.8
· 60	488	340	317	1,456	274	63,940	<17	<6.8
61 -	545	<243	317		<211	47,060	<17	<6.8
0 0	<287	292	1,561	3,587	1,625	141,600	<17	×6.8
) [<287	<243	<211	401	<211	12,000	<17	<6.8
	<287	<243	<211	422	<211	14,390	<17	<6.8

TCE = TRICHLOROETHENE PCE = TETRACHLOROETHENE

^{*} BIEX WERE ANALYZED VIA GC/FID AND HALOCARBONS WERE ANALYZED VIA GC/ECD.

^{&#}x27;CALCULATED USING THE SUM OF THE AREAS OF ALL INTEGRATED CHROMATOGRAM PEAKS AND THE INSTRUMENT RESPONSE FACTOR FOR TOLUENE

TABLE 2 (cont)

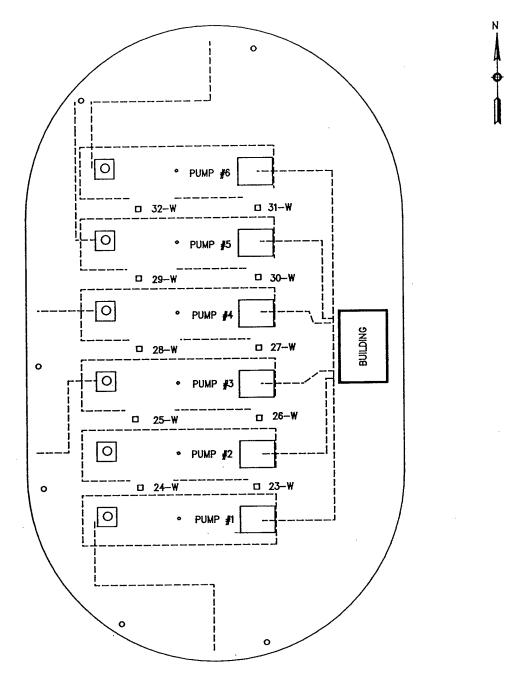
* LABORATORY RESULTS OF SOIL GAS SAMPLES CONCENTRATIONS IN PARTS PER BILLION (V/V)

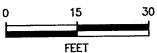
SAMPLE	BENZENE	TOLUENE	ETHYL- BENZENE	m- & p- XYLENE	O- XYLENE	TOTAL VOLATILES1	TCE	PCE
FIELD CONTROL SAMPLES	OL SAMPLES	zol.	÷*					
, H	<287	<243	<211	<211	<211	<243	<17	<6.8
	<287	<243	<211	<211	<211	<243	<17	8.9>
i, m	<287	<243	<211	<211	<211	<243	<17	×6.8
7	<287	<243	<211	<211	<211	<243	<17	×6.8
LABORATORY DUPLICATE ANALYSES	DUPLICATE	ANALYSES						
16	<287	<243	<211	359	<211	7,333	<17	<6.8
16R	<287	<243	253	359	<211	7,302	<17	<6.8
LABORATORY BLANKS	BLANKS		:		•			
BMCHM-1	<287	<243	<211	<211	<211	<243	<17	<6.8
							,	

^{*} BTEX WERE ANALYZED VIA GC/FID AND HALOCARBONS WERE ANALYZED VIA GC/ECD.

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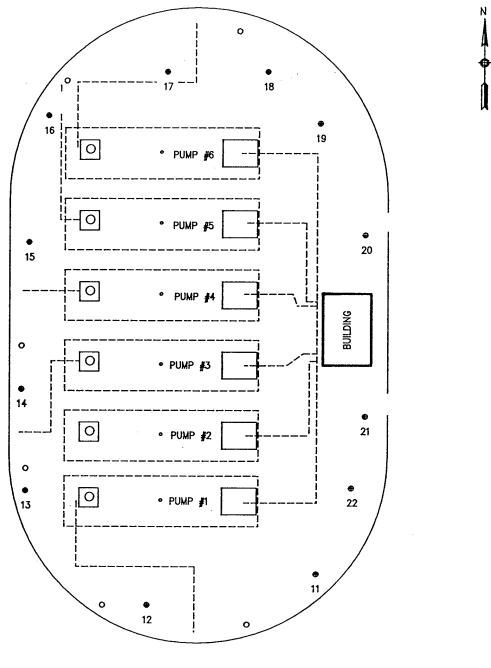
- WATER SAMPLE LOCATION
- o TANK FILL

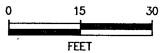
FIGURE 1A. Ground Water Sample Locations



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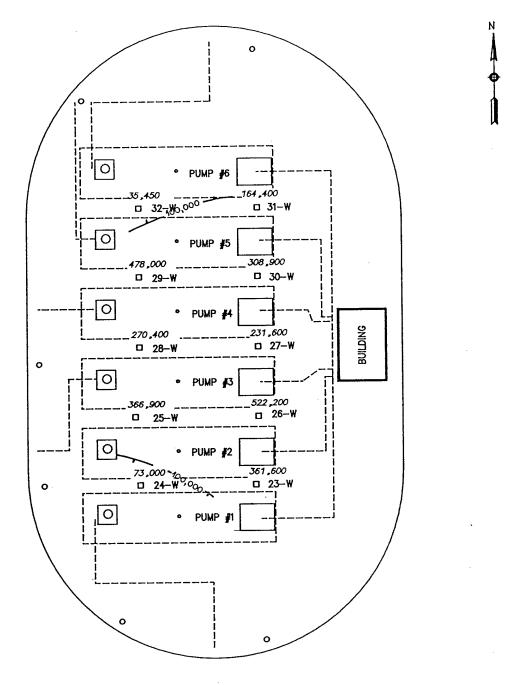
- SOIL GAS SAMPLE LOCATION
- o TANK FILL

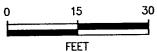
FIGURE 1B. Soil Gas Sample Locations



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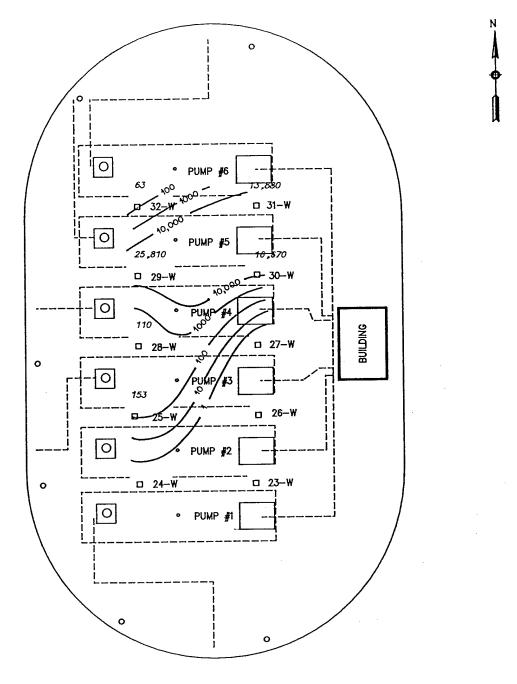


- **D** WATER SAMPLE LOCATION
- O TANK FILL



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FIGURE 2. FID Total Volatiles in Ground Water (calc'd ppb v/v)



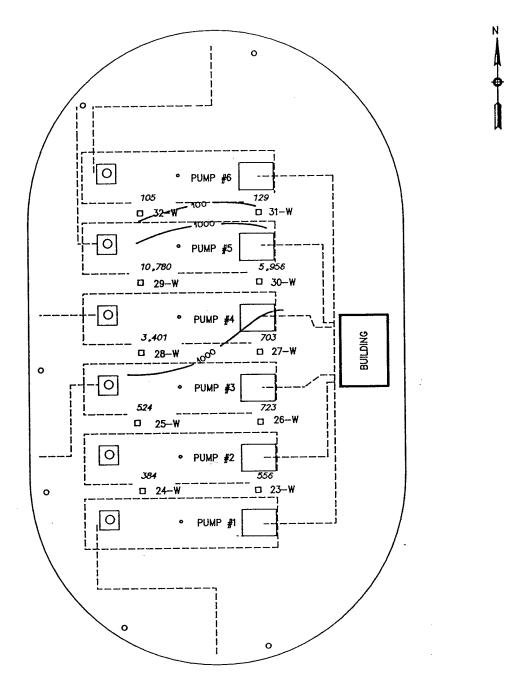


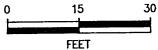
- D WATER SAMPLE LOCATION
- o TANK FILL

FIGURE 3. Benzene in Ground Water (ppb v/v)



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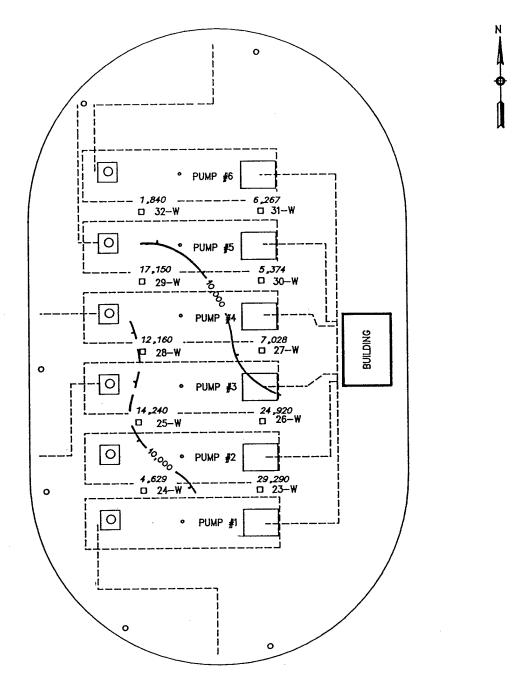


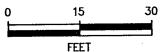
- D WATER SAMPLE LOCATION
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FIGURE 4. Toluene in Ground Water (ppb v/v)



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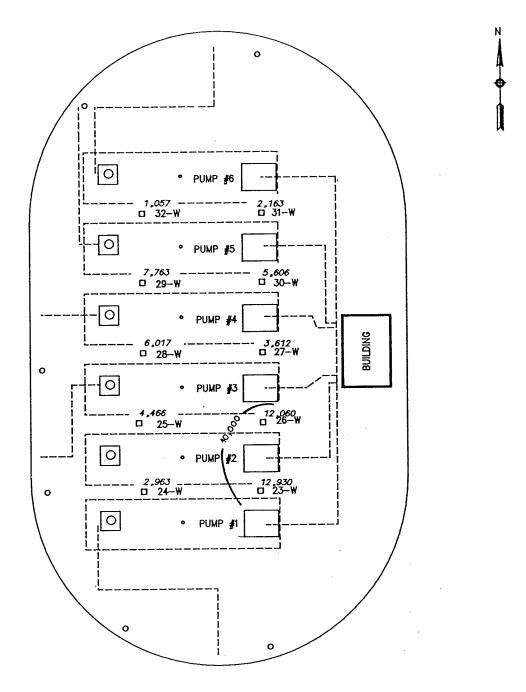
- WATER SAMPLE LOCATION
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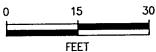
FIGURE 5. meta— and para—Xylene in Ground Water (ppb v/v)



ENVIRONMENTAL SERVICES, INC.

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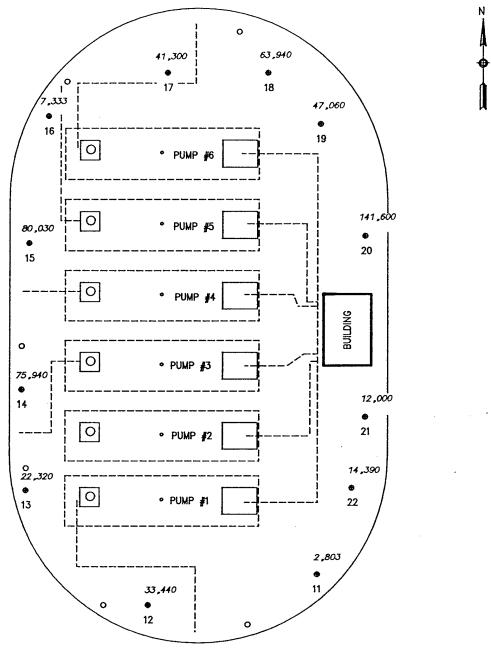


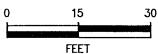
- □ WATER SAMPLE LOCATION
- o TANK FILL



This map is integral to a written report and should be viewed in that context.

FIGURE 6. ortho—Xylene in Ground Water (ppb v/v)





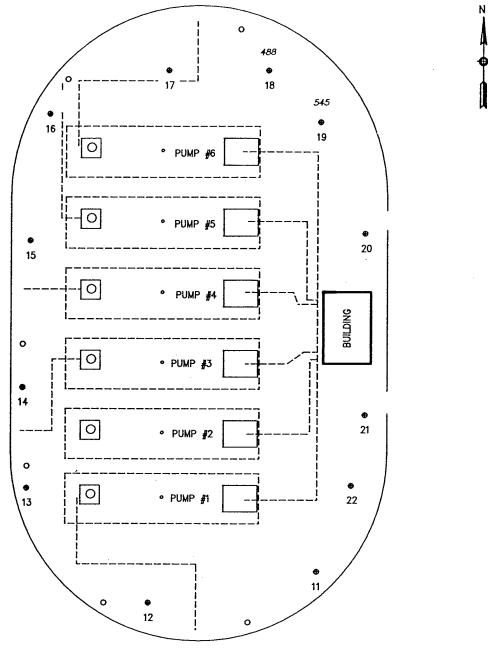
- SOIL GAS SAMPLE LOCATION
- o TANK FILL

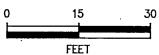
FIGURE 7. FID Total Volatiles in Soil Gas (calc'd ppb v/v)



ENVIRONMENTAL SERVICES, INC.

This map is integral to a written report and should be viewed in that context.



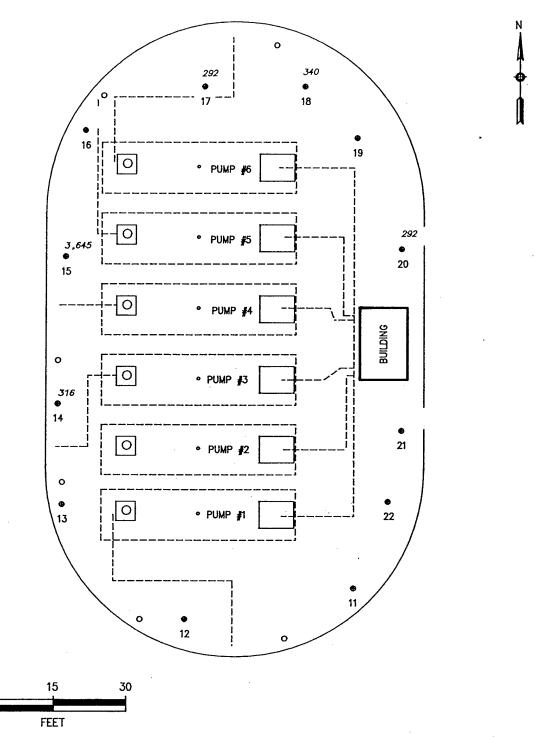


- SOIL GAS SAMPLE LOCATION
- o TANK FILL

FIGURE 8. Benzene in Soil Gas (ppb)



This map is integral to a written report and should be viewed in that context.



• SOIL GAS SAMPLE LOCATION

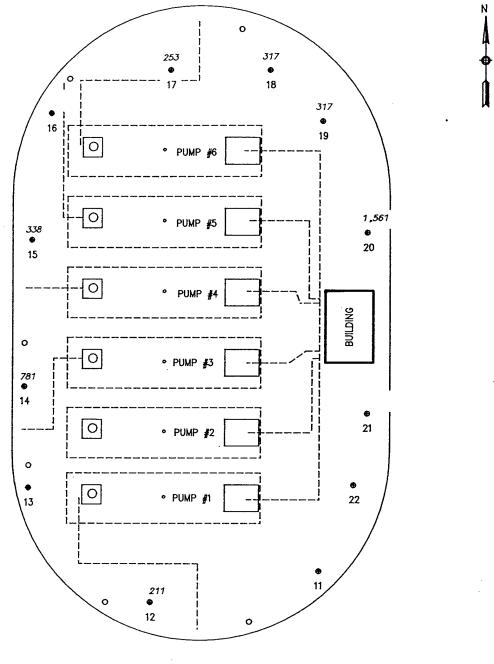
o TANK FILL

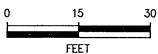
FIGURE 9. Toluene in Soil Gas (ppb)



ENVIRONMENTAL SERVICES, INC.

This map is integral to a written report and should be viewed in that context.



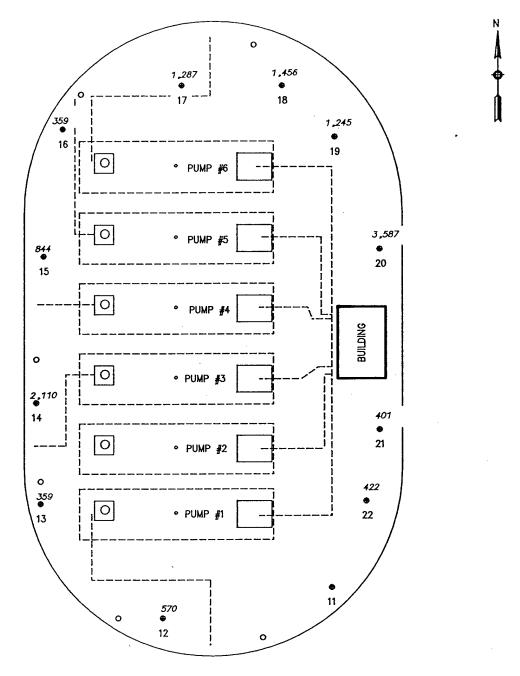


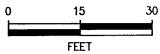
- SOIL GAS SAMPLE LOCATION
- o TANK FILL -

FIGURE 10. Ethylbenzene in Soil Gas (ppb)



This map is integral to a written report and should be viewed in that context.



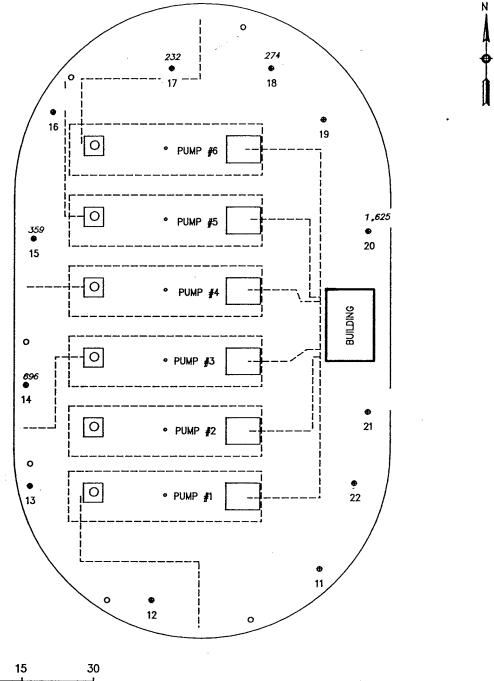


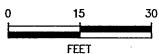
- SOIL GAS SAMPLE LOCATION
- o TANK FILL

FIGURE 11. meta— and para—Xylene in Soil Gas (ppb)



This map is integral to a written report and should be viewed in that context.





- SOIL GAS SAMPLE LOCATION
- o TANK FILL

FIGURE 12. ortho—Xylene in Soil Gas (ppb)



ENVIRONMENTAL SERVICES, INC.

This map is integral to a written report and should be viewed in that context.

GEOTECHNICAL RESULTS

Box Sample (no label)

3516 Greensboro Avenue • P.O. Drawer 1128 • Tuscaloosa, Alabama 35403 • Telephone 205-345-0816 • FAX 205-345-0992

SOIL ANALYSES REPORT

KILMAN BROTHERS, INC. DANNELLY A.N.G. FIELD MGM27526.S.I.M.G. MAY, 1991

SAMPLE I.D.	DEPTH	PERMEABILITY COEFFICIENT "k"	INSITU DRY UNIT WEIGHT	AS RECEIVED MOISTURE
OWS-3 (Site 2)	8'-10'	2.3 x 10 ⁻⁶ cm/sec	89.7 pcf	31.5 %
OWS-4 (Site 2)	8'-10'	1.3 x 10 ⁻⁸ cm/sec	91.6 pcf	29.0 %
POL-8	10'-12'	7.2 x 10 ⁻⁹ cm/sec	88.3 pcf	34.7 %
•	٠.			
		CALCIUM CARBONAT	E DATA	
	SAMPLE TIFICATION	CALCI	UM CARBONATE (% CaCO ₃)	EQUIVALENT

TTL, Inc.

Britette L. Lee, Geologist

35.6 %

SUMMARY TABLES OF ANALYTICAL DATA

DANNELLY ANG - Montgomery, Alabama (*) BACKGROUND INFORMATION 187th Fighter Group

Client Somnle ID				B1P-S-0-2	B1P-S-8.5-10	B1P-S-18.5-20	BIP-S-0-2 BIP-S-8.5-10 BIP-S-18.5-20 BIP-S18.5-20D B3P-S-8.5-10 B3P-S-13.5-15 B3P-S-28.5-30	B3P-S-8.5-10	B3P-S-13.5-15	B3P-S-28.5-30
Lab Sample Number				28909001	28909002	28909003	28909004	28922001	28922002	28922003
Matrix				SOIL	ł	SOIL	SOIL	TIOS	SOIL	SOIL
Sample Date				2/26/91	2/26/91	2/26/91	2/26/91	2/27/91	2/27/91	2/27/91
	MCL	CRDL	Action Levels							
Volatile Compounds	ng/L	ug/kg	ug/kg/ug/L				· · · · · · · · · · · · · · · · · · ·			
2-Butanone	N/A	10	N/A/N/A							
Acetone	N/A	10	8E+06/4,000	16 B	33 B	110B	55 B	13 B	15 B	31B
Carbon Disulfide	N/A	5	8E+06/4,000							
Ethyl Benzene	7E+02	5	8E+06/4,000							6 J
Methylene Chloride	N/A	5	5/000/6	8 B	31 B	35 B	33 B	8 B	7 B	8 B
Toluene	1E+03	5	2E+08/10,000					1 J		3 J
Xvlene (total)	1E+04	5	2E+08/70,000							4 J
Chloromethane	N/A	10	N/A/N/A							
Bromomethane	N/A	10	100,000/N/A							
Vinyl Chloride	2	10	N/A/N/A							
Chloroethane	N/A	10	N/A/N/A							
1,1-Dichloroethene	N/A	5	N/A/N/A							
Acrylonitrile	N/A		1000/0.06							

B - Applies to organic data only. Present in the corresponding method blank.

J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.

MCL - Maximum Contaminant Level (U.S. Drinking Water Standards)

CRDL - Contract Required Detection Limit

Action levels proposed in Appendix A of 40CFR254.521(a) Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

(a) - Data reported in ug/kg (soil) and ug/L (water).

DANNELLY ANG - Montgomery, Alabama (*)

Client Sample ID				B1P-S-0-2	B1P-S-8.5-10	B1P-S-18.5-20	BIP-S-0-2 BIP-S-8.5-10 BIP-S-18.5-20 BIP-S18.5-20D B3P-S-8.5-10 B3P-S-13.5-15 B3P-S-28.5-30	B3P-S-8.5-10	B3P-S-13.5-15	B3P-S-28.5-30
Lab Sample Number				28909001	28909001 28909002	28909003	28909004	28922001	28922002	28922003
Matrix				SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL
Sample Date				2/26/91	2/26/91	16/97/2	16/97/2	16/12/7	2/27/91	2/27/91
	MCL	CRDL	Action Level							
Volatile Compounds	ng/L	ug/kg	ug/kg/ug/L							
1,1-Dichloroethane	N/A	5	N/A/N/A							
1,2-Dichloroethene (total)	N/A	5	8,000/N/A							
Chloroform	1E+02	5	N/A/N/A							
1,2-Dichloroethane	5	5	20,000/400							
1,1,1-Trichloroethane	2E+02	5	2E+06/700							
Carbon Tetrachloride	5	5	N/A/N/A			-				
Vinyl Acetate	N/A	10	N/A/N/A							
Bromodichloromethane	1E+02	5	500/0.02							
1,2-Dichloropropane	5	5	N/A/N/A							
cis-1,3-Dichloropropene	N/A	5	40,000/20							
Trichloroethene	5	5	2E+06/700							
Dibromochloromethane	N/A	5	2E+06/7,000							

Notes:

B - Applies to organic data only. Present in the corresponding method blank.

J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.

MCL - Maximum Contaminant Level (U.S. Drinking Water Standards)

CRDL - Contract Required Detection Limit

Action levels proposed in Appendix A of 40CFR254.521(a)

Blank Space - Value is below the Method Detection Limit (MDL).

N/A - Not Applicable

(a) - Data reported in ug/kg (soil) and ug/L (water).

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Client Sample ID				B1P-S-0-2	B1P-S-8.5-10	B1P-S-18.5-20	BIP-S18.5-20D	B3P-S-8.5-10	B3P-S-13.5-15	B3P-S-28.5-30
Lab Sample Number				28909001	28909002	28909003	28909004	28922001	28922002	28922003
Matrix				SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL
Sample Date				2/26/91	2/26/91	2/26/91	2/26/91	2/27/91	2/27/91	2/27/91
· 第二十八十二十二十二十二十二十二十二十二十二十二十二十二十二十二十二十二十二十二	MCL	CRDL	Action Levels						を受けるというと言葉を	では、大学の学の意味を
Volatile Compounds	ng/L	ug/kg	ug/kg/ug/L				三十二十二十二十二十二十二十二十二十二十二十二十二十二十二十二十二十二十二十二			
1,1,2-Trichloroethane	5	5	100,000/6.0							
Benzene	2	5	N/A/N/A							
trans-1,3-Dichloropropene	N/A	5	20,000/400							
Bromoform	1E+02	5	2E+06/700							
4-Methyl-2-Pentanone	A/N	10	N/A/N/A							
2-Hexanone	A/A	10	N/A/N/A							
Tetrachloroethene	2	5	N/A/N/A							
1,1,2,2-Tetrachloroethane	N/A	5	40,000/20							
Chlorobenzene	N/A	5	2E+06/700							
Styrene	1E+02	5	2E+06/7,000							

B - Applies to organic data only. Present in the corresponding method blank.

J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.

MCL - Maximum Contaminant Level (U.S. Drinking Water Standards)

CRDL - Contract Required Detection Limit

Action levels proposed in Appendix A of 40CFR254.521(a) Blank Space - Value is below the Method Detection Limit (MDL).

N/A - Not Applicable (a) - Data reported in ug/kg (soil) and ug/L (water).

DANNELLY ANG - Montgomery, Alabama (*)

Client Sample ID					B5P-S-0-2	B5P-S-8.5-10	B5P-S-23.0-25	B5P-S-0-2 B5P-S-8.5-10 B5P-S-23.0-25 B5P-S-23.0-25D BG-MW-6	BG-MW-6	BG-2P
Lab Sample Number					28909005	28909006	28909007	28909008	18301001	18311001
Matrix					SOIL	SOIL	SOIL	SOIL	WATER	WATER
Sample Date					2/26/91	2/26/91	2/26/91	2/26/91	16/01/4	4/12/91
を書きたける。 は、これでは、一般では、これでは、これでは、これでは、これでは、これでは、これでは、これでは、これ	MCL	CRDL	CRDL	Action Levels						
Volatile Compounds	ug/L	ug/kg	ug/L	ug/kg/ug/L						
2-Butanone	N/A	10	10	N/A/N/A			8 J			
Acetone	N/A	10	10	8E+06/4,000	27 B	78 B	110 B		16 B	
Carbon Disulfide	N/A	5	5	8E+06/4,000			10			
Ethyl Benzene	7E+02	S	5	8E+06/4,000						
Methylene Chloride	N/A	5	5	5/000'06	30 B	9 B	11 B		3 B	
Toluene	1E+03	S	5	2E+08/10,000						
Xylene (total)	1E+04	5	5	2E+08/70,000					1B	
Chloromethane	N/A	10	10	N/A/N/A						
Bromomethane	N/A	10	10	100,000/N/A						
Vinyl Chloride	2	10	10	N/A/N/A						
Chloroethane	N/A	10	10	N/A/N/A			•			
1,1-Dichloroethene	N/A	5	5	N/A/N/A			,			
Acrylonitrile	N/A			1,000/0.06						100 B

Notes:

B - Applies to organic data only. Present in the corresponding method blank.

J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.

MCL - Maximum Contaminant Level (U.S. Drinking Water Standards)

CRDL - Contract Required Detection Limit

Action levels proposed in Appendix A of 40CFR254.521(a) Blank Space - Value is below the Method Detection Limit(MDL). N/A - Not Applicable (a) - Data reported in ug/kg (soil) and ug/L (water).

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DANNELLY ANG - Montgomery, Alabama (a) BACKGROUND INFORMATION 187th Fighter Group

Client Sample ID					B5P-S-0-2	B5P-S-8.5-10	B5P-S-23.0-25	B5P-S-0-2 B5P-S-8.5-10 B5P-S-23.0-25 B5P-S-23.0-25D BG-MW-6	BG-MW-6	BG-2P
Lab Sample Number					28909005	28909005 28909006	28909007	28909008	18301001	18311001
Matrix		:			SOIL	TIOS	SOIL	TIOS	WATER	WATER
Sample Date					2/26/91	16/97/7	2/26/91	16/97/7	4/10/91	4/12/91
松本の日本の日本の日本の日本の日本の日本の日本の日本の日本の日本の日本の日本の日本	MCL	CRDL	CRDL	Action Levels						the state of the s
Volatile Compounds	ug/L	ug/kg	ng/L	ug/kg/ug/L						
1,1-Dichloroethane	N/A	5	5	N/A/N/A						
1,2-Dichloroethene (total)	N/A	5	5	8,000/N/A						
Chloroform	1E+02	5	5	100,000/6.0						
1,2-Dichlorocthane	5	5	5	N/A/5.0						
1,1,1-Trichloroethane	2E+02	5	5	7E+06/3,000						
Carbon Tetrachloride	5	5	5	5000/0.3						
Vinyl Acetate	N/A	10	10	N/A/N/A						
Bromodichloromethane	1E+02	5	5	500/0.02						
1,2-Dichloropropane	5	5	5	N/A/N/A						
cis-1,3-Dichloropropene	N/A	5	5	20,000/10.0						
Trichloroethene	5	5	5	N/A/N/A						
Dibromochloromethane	N/A	5	5	N/A/N/A						

B - Applies to organic data only. Present in the corresponding method blank.

J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.

MCL - Maximum Contaminant Level (U.S. Drinking Water Standards)

CRDL - Contract Required Detection Limit

Action levels proposed in Appendix A of 40CFR254.521(a) Blank Space - Value is below the Method Detection Limit(MDL). N/A - Not Applicable (a) - Data reported in ug/kg (soil) and ug/L (water).

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DANNELLY ANG - Montgomery, Alabama (a)

Client Sample ID					B5P-S-0-2	B5P-S-8.5-10	B5P-S-23.0-25	B5P-S-0-2 B5P-S-8.5-10 B5P-S-23.0-25 B5P-S-23.0-25D BG-MW-6 BG-2P	BG-MW-6	BG-2P
Lab Sample Number					28909005	28909005 28909006	28909007	28909008	18301001	18311001
Matrix					SOIL	SOIL	SOIL	SOIL	WATER	WATER
Sample Date					2/26/91	2/26/91	2/26/91	2/26/91	4/10/91	4/12/91
	MCL	CRDL	CRDL	Action Levels						工艺學學學
Volatile Compounds	ng/L	ug/kg	ug/L	ug/kg/ug/L		地名加拿加加纳				
trans-1,3-Dichloropropene	N/A	5	5	20,000/400						
Bromoform	1E+02	5	5	2E+06/700						
4-Methyl-2-Pentanone	N/A	10	10	N/A/N/A						
2-Hexanone	N/A	10	10	N/A/N/A						
Tetrachloroethene	5	5	5	N/A/N/A						
1,1,2,2-Tetrachloroethane	N/A	5	5	40,000/20						
Chlorobenzene	N/A	5	5	2E+06/700						
Styrene	1E+02	5	5	2E+06/7,000						

B - Applies to organic data only. Present in the corresponding method blank.

J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.

MCL - Maximum Contaminant Level (U.S. Drinking Water Standards)

CRDL - Contract Required Detection Limit

Action levels proposed in Appendix A of 40CFR254.521(a) Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

(a) - Data reported in ug/kg (soil) and ug/L (water).

DANNELLY ANG - Montgomery, Alabama "

Transfer of			B1P-S-0-2	B1P-S-8.5-10	B1P-S-18.5-20	B1P-S18.5-20D	B3P-S-8.5-10	B3P-S-13.5-15	B3P-S-28.5-30
Tob Sample Mumber			28909001	28909002	28909003	i	28922001	28922002	28922003
Mateix			SOIL	SOIL	SOIL	SOIL	SOIL	TIOS	SOIL
Comple Date			2/26/91	2/26/91	2/26/91	2/26/91	2/27/91	2/27/91	2/27/91
	CRDL	Action Levels						から 一年の後の一年の日本の	がおける (And And And And And And And And And And
Semivolatile Compounds	ug/kg	ug/kg		建筑的新疆域的					
Fluoranthene	330	A/N	79.1						
Phenanthrene	330	N/A	52 J						
Pyrene	330	N/A	56 J						
his(2-Ethylhexyl)-phthalate	330	20,000		1600 B	1700 B		730 B	1200 B	1100 B
Phenol	330	5E+07							
bis(2-Chloroethyl)Ether	330	009							
2-Chlorophenol	330	400,000							
1,3-Dichlorobenzene	330	N/A							
1,4-Dichlorobenzene	330	N/A	1						
Benzyl Alcohol	330	N/A							
1,2-Dichlorobenzene	330	N/A	-						
2-Methylphenol	330	N/A							

B - Applies to organic data only. Present in the corresponding method blank.
 J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.
 CRDL - Contract Required Detection Limit

Action levels proposed in Appendix A of 40CFR254.521(a) Blank Space - Value is below the Method Detection Limit(MDL). N/A - Not Applicable (a) - Data reported in ug/kg.

BACKGROUND INFORMATION

DANNELLY ANG - Montgomery, Alabama (*) 187th Fighter Group

Client Sample ID			BIP-S-0-2	B1P-S-8.5-10	BIP-S-18.5-20	B1P-S-8.5-10 B1P-S-18.5-20 B1P-S18.5-20D B3P-S-8.5-10 B3P-S-13.5-15 B3P-S-28.5-30	B3P-S-8.5-10	B3P-S-13.5-15	B3P-S-28.5-30
Lab Sample Number			28909001	28909002	28909003	28909004	28922001	28922002	28922003
Matrix			SOIL	TIOS	SOIL	SOIL	SOIL	SOIL	SOIL
Sample Date			2/26/91	2/26/91	2/26/91	2/26/91	2/27/91	2/27/91	2/27/91
	CRDL	Action Levels						主动之际重要重要	
Semivolatile Compounds	ug/kg	ug/kg				大の子が、これをおり、日本の日本の日本の日本の日本の日本の日本の日本の日本の日本の日本の日本の日本の日	电影图像影响影响唱		
ois(2-Chloroisopropyl)Ether	330	ΝΆ							
f-Methylphenol	330	N/A							
N-Nitroso-di-n-propylamine	330	100							
fexachloroethane	330	80,000							
Vitrobenzene	330	40,000							
sophorone	330	2E+06							
2-Nitrophenol	330	N/A							
2,4-Dimethylphenol	330	N/A							
3enzoic Acid	1600	N/A							
ois(2-Chloroethoxy)Methane	330	N/A							
2,4-Dichlorophenol	330	200,000							
1,2,4-Trichlorobenzene	330	2E+06							

Notes:

B - Applies to organic data only. Present in the corresponding method blank.
J - Applies to organic data only. Value detected is greater than zero but less than the CRDL. CRDL - Contract Required Detection Limit
Action levels proposed in Appendix A of 40CFR254.521(a)
Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable (a) - Data reported in ug/kg.

DANNELLY ANG - Montgomery, Alabama "

Client Somule ID			B1P-S-0-2	B1P-S-8.5-10	B1P-S-18.5-20	B1P-S18.5-20D	B3P-S-8.5-10	B3P-S-13,5-15 B3P-S-28,5-30	B3P-S-28.5-30
I ah Sample Number			28909001	28909002	28909003	28909004	28922001	28922002	28922003
Matrix			SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL
Sample Date			2/26/91	2/26/91	2/26/91	2/26/91	2/27/91	2/27/91	2/27/91
Sample Date	CRDL	Action Levels						"妈妈妈"	
Semivolatile Compounds	ug/kg					3 2 3 3 3 3 3 3 3 3 3 3 3 3 3 3 3 3 3 3			
Naphthalene	330	N/A							
4-Chloroaniline	330	N/A							
Hexachlorobutadiene	330	000'06							
4-Chloro-3-methylphenol	330	N/A							
2-Methylnaphthalene	330	N/A							
Hexachlorocyclopentadiene	330	000,009							
2.4.6-Trichlorophenol	330	40,000							
2,4,5-Trichlorophenol	1600	8E+06							
2-Chloronaphthalene	330	N/A							
2-Nitroaniline	1600	N/A							
Dimethyl Phthalate	330	N/A							
Acenaphthylene	330	N/A							

B - Applies to organic data only. Present in the corresponding method blank.
 J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.
 CRDL - Contract Required Detection Limit

Action levels proposed in Appendix A of 40CFR254.521(a) Blank Space - Value is below the Method Detection Limit(MDL). N/A - Not Applicable (a) - Data reported in ug/kg.

DANNELLY ANG - Montgomery, Alabama (*) BACKGROUND INFORMATION 187th Fighter Group

Client Sample ID			B1P-S-0-2	B1P-S-8.5-10	B1P-S-18.5-20	BIP-S18.5-20D	B3P-S-8.5-10	B3P-S-13.5-15	B3P-S-28.5-30
Lab Sample Number			28909001	28909002	28909003	28909004	28922001	28922002	28922003
Matrix			TIOS	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL
Sample Date			16/97/7	2/26/91	2/26/91	2/26/91	2/27/91	2/27/91	2/27/91
	CRDL	Action Levels		學者 医神经病 医三角					
Semivolatile Compounds	ug/kg	ug/kg							
2,6-Dinitrotoluene	330	1000							
3-Nitroaniline	1600	N/A							
Acenaphthene	330	N/A							
2,4-Dintrophenol	1600	200,000							
4-Nitrophenol	1600	N/A							
Dibenzofuran	330	N/A							
2,4-Dinitrotoluene	330	N/A							
Diethylphthalate	330	6E+07							
4-Chlorophenyl-phenylether	330	N/A							-
Fluorene	330	N/A							
4-Nitroaniline	1600	N/A							
4,6-Dinitro-2-methylphenol	1600	N/A							

B - Applies to organic data only. Present in the corresponding method blank.
 J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.
 CRDL - Contract Required Detection Limit

Action levels proposed in Appendix A of 40CFR254.521(a)

Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable (a) - Data reported in ug/kg.

DANNELLY ANG - Montgomery, Alabama (*)

Client Sample ID			B1P-S-0-2	B1P-S-8.5-10	B1P-S-18.5-20	B1P-S18.5-20D	B3P-S-8.5-10	B3P-S-13.5-15	B3P-S-28.5-30
Lab Sample Number			28909001	28909002	28909003	28909004	28922001	28922002	28922003
Matrix			TIOS	TIOS	TIOS	SOIL	SOIL	SOIL	SOIL
Sample Date			2/26/91	2/26/91	2/26/91	2/26/91	2/27/91	2/27/91	2/27/91
このない 一次のでは、一般の一般の一般の一般の一般の一般の一般の一般の一般の一般の一般の一般の一般の一	CRDL	Action Levels							
Semivolatile Compounds	ug/kg	ug/kg							
N-Nitrosodiphenylamine(1)	330	100,000							
4-Bromophenyl-phenylether	330	N/A							
Hexachlorobenzene	330	N/A							
Pentachiorophenol	1600	2E+06		,					
Anthracene	330	N/A							
Di-n-Buthylphthalate	330	8E+06							
Butylbenzylphthalate	330	2E+07							
3,3'-Dichlorobenzidine	099	2000							
Benzo(a)anthracene	330	N/A							
Chrysene	330	N/A							
Di-n-octylphthalate	330	N/A			-				
Benzo(b)fluoranthene	330	N/A							

Notes:

B - Applies to organic data only. Present in the corresponding method blank. J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.

CRDL - Contract Required Detection Limit

Action levels proposed in Appendix A of 40CFR254.521(a) Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable (a) - Data reported in ug/kg.

DANNELLY ANG - Montgomery, Alabama (a)

Client Sample ID			B1P-S-0-2	B1P-S-8.5-10	B1P-S-18.5-20	B1P-S18.5-20D	B3P-S-8.5-10	B3P-S-13.5-15	B3P-S-28,5-30
Lab Sample Number			28909001	28909002	28909003	28909004	28922001	28922002	28922003
Matrix			TIOS	TIOS	SOIL	TIOS	SOIL	SOIL	SOIL
Sample Date			2/26/91	2/26/91	2/26/91	2/26/91	2/27/91	2/27/91	2/27/91
	CRDL	Action Levels							
Semivolatile Compounds	ug/kg	ug/kg	は最高をできない。			· · · · · · · · · · · · · · · · · · ·	以上の表現の表現の表現である。 のは、 のは、 のは、 のは、 のは、 のは、 のは、 のは、		
Benzo(b)fluoranthene	330	N/A							
Benzo(k)fluoranthene	330	N/A							
Benzo(a)pyrene	330	N/A							
Indeno(1,2,3-cd)Pyrene	330	N/A							
Dibenz(a,h)Anthracene	330	N/A							
benzo(g,h,i)perylene	330	N/A							

Notes:

B - Applies to organic data only. Present in the corresponding method blank.

J - Applies to organic data only. Value detected is greater than zero but less than the CRDL. CRDL - Contract Required Detection Limit

Action levels proposed in Appendix A of 40CFR254.521(a) Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable (a) - Data reported in ug/kg.

C-12

BACKGROUND INFORMATION 187th Fighter Group DANNELLY ANG - Montgomery, Alabama (*)

Client Sample ID			B5P-S-0-2	B5P-S-8.5-10	B5P-S-23.0-25	B5P-S-23.0-25D
Lab Sample Number			28909005	28909006	28909007	28909008
Matrix			SOIL	SOIL	TIOS	SOIL
Sample Date			2/26/91	2/26/91	2/26/91	2/26/91
	CRDL	Action Levels				· · · · · · · · · · · · · · · · · · ·
Semivolatile Compounds	ug/kg	ug/kg				
Fluoranthene	330	N/A				
Phenanthrene	330	N/A				
Pyrene	330	N/A				
bis(2-Ethylhexyl)-phthalate	330	50,000	970 B	2500 B	3700 B	1400 B
Phenol	330	5E+07				
bis(2-Chloroethyl)Ether	330	009				
2-Chlorophenol	330	400,000				
1,3-Dichlorobenzene	330	N/A				
1,4-Dichlorobenzene	330	N/A				
Benzyl Alcohol	330	N/A				
1,2-Dichlorobenzene	330	N/A				
2-Methylphenol	330	N/A				

Notes

B - Applies to organic data only. Present in the corresponding method blank.

J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.

CRDL - Contract Required Detection Limit

Action levels proposed in Appendix A of 40CFR254.521(a)

Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

(a) - Data reported in ug/kg.

C-13

DANNELLY ANG - Montgomery, Alabama (*)

Client Sample ID			B5P-S-0-2	B5P-S-8.5-10	B5P-S-23.0-25	B5P-S-23.0-25D
Lab Sample Number			28909005	28909006	28909007	28909008
Matrix	٠		SOIL	SOIL	SOIL	SOIL
Sample Date		-	2/26/91	2/26/91	2/26/91	2/26/91
	CRDL	Action Levels				
Semivolatile Compounds	ug/kg	ug/kg				
bis(2-Chloroisopropyl)Ether	330	N/A				
4-Methylphenol	330	N/A				
N-Nitroso-di-n-propylamine	330	100				
Hexachloroethane	330	80,000				
Nitrobenzene	330	40,000				
Isophorone	330	2E+06				
2-Nitrophenol	330	N/A	-			
2,4-Dimethylphenol	330	N/A				
Benzoic Acid	1600	N/A				
bis(2-Chloroethoxy)Methane	330	N/A				
2,4-Dichlorophenol	330	200,000				
1,2,4-Trichlorobenzene	330	2E+06				

Notes:

B - Applies to organic data only. Present in the corresponding method blank.

J - Applies to organic data only. Value detected is greater than zero but less than the CRDL. CRDL - Contract Required Detection Limit

Action levels proposed in Appendix A of 40CFR254.521(a)

Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

DANNELLY ANG - Montgomery, Alabama (a)

Client Sample ID			B5P-S-0-2	B5P-S-8.5-10	B5P-S-23.0-25	B5P-S-8.5-10 B5P-S-23.0-25 B5P-S-23.0-25D
Lab Sample Number			28909005	28909006	28909007	28909008
Matrix			SOIL	SOIL	SOIL	SOIL
Sample Date		-	2/26/91	2/26/91	2/26/91	2/26/91
	CRDL	Action Levels				
Semivolatile Compounds	ug/kg	ug/kg				
Naphthalene	330	N/A				
4-Chloroaniline	330	N/A				
Hexachlorobutadiene	330	000'06				
4-Chloro-3-methylphenol	330	N/A				
2-Methylnaphthalene	330	N/A				
Hexachlorocyclopentadiene	330	000,009				
2,4,6-Trichlorophenol	330	40,000				
2,4,5-Trichlorophenol	1600	8E+06				
2-Chloronaphthalene	330	N/A				
2-Nitroaniline	1600	N/A				
Dimethyl Phthalate	330	N/A				
Acenaphthylene	330	N/A				

Notes:

B - Applies to organic data only. Present in the corresponding method blank.

J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.

CRDL - Contract Required Detection Limit

Action levels proposed in Appendix A of 40CFR254.521(a)

Blank Space - Value is below the Method Detection Limit(MDL). N/A - Not Applicable

DANNELLY ANG - Montgomery, Alabama (*)

an classical survivor			טטטטט	DED 0 0 5 10	D C D C D D C D C	D5D 6 73 0 75D
Cilent Sample 1D			B3F-3-0-2	B3F-3-6.3-10	BJF-5-25.0-23	D2F-3-43.0-43D
Lab Sample Number			28909005	90060687	28909007	28909008
Matrix			SOIL	TIOS	TIOS	SOIL
Sample Date			2/26/91	2/26/91	2/26/91	2/26/91
	CRDL	Action Levels				
Semivolatile Compounds	ug/kg	ug/kg				
2,6-Dinitrotoluene	330	1000				
3-Nitroaniline	1600	N/A				
Acenaphthene	330	N/A				
2,4-Dintrophenol	1600	200,000				
4-Nitrophenol	1600	N/A				
Dibenzofuran	330	N/A				
2,4-Dinitrotoluene	330	N/A				
Diethylphthalate	330	6E+07				
4-Chlorophenyl-phenylether	330	N/A				
Fluorene	330	N/A				
4-Nitroaniline	1600	N/A				
4,6-Dinitro-2-methylphenol	1600	N/A				

Notes

B - Applies to organic data only. Present in the corresponding method blank.

J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.

CRDL - Contract Required Detection Limit

Action levels proposed in Appendix A of 40CFR254.521(a) Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

DANNELLY ANG - Montgomery, Alabama (*)

Client Sample ID			B5P-S-0-2	B5P-S-8.5-10	B5P-S-23.0-25	B5P-S-23.0-25D
Lab Sample Number			28909005	28909006	28909007	28909008
Matrix			SOIL	SOIL	SOIL	SOIL
Sample Date	CRDL	Action Levels	2/26/91	2/26/91	2/26/91	2/26/91
大学 のから かんかん かんかん かんかん かんかん かんかん かんかん かんかん	ug/kg	ug/kg				
N-Nitrosodiphenylamine(1)	330	100,000				
4-Bromophenyl-phenylether	330	N/A				
Hexachlorobenzene	330	N/A				
Pentachlorophenol	1600	2E+06				
Anthracene	330	N/A				
Di-n-Buthylphthalate	330	8E+06				
Butylbenzylphthalate	330	2E+07				
3,3'-Dichlorobenzidine	099	2000				
Benzo(a)anthracene	330	N/A				
Chrysene	330	N/A				
Di-n-octylphthalate	330	N/A				
Benzo(b)fluoranthene	330	N/A				

B - Applies to organic data only. Present in the corresponding method blank.

J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.

CRDL - Contract Required Detection Limit

Action levels proposed in Appendix A of 40CFR254.521(a) Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

DANNELLY ANG - Montgomery, Alabama (*)

Client Sample ID			B5P-S-0-2	B5P-S-8.5-10		B5P-S-23.0-25 B5P-S-23.0-25D
Lab Sample Number			28909005	28909006		28909008
Matrix		-	TIOS	SOIL		SOIL
Sample Date		-	2/26/91	2/26/91	2/26/91	2/26/91
	CRDL	Action Levels				
Semivolatile Compounds	ug/kg	ug/kg				
Benzo(k)fluoranthene	330	N/A				
Benzo(a)pyrene	330	N/A				
Indeno(1,2,3-cd)Pyrene	330	N/A				
Dibenz(a,h)Anthracene	330	N/A			-	
benzo(g,h,i)perylene	330	N/A				

B - Applies to organic data only. Present in the corresponding method blank.

J - Applies to organic data only. Value detected is greater than zero but less than the CRDL. CRDL - Contract Required Detection Limit

Action levels proposed in Appendix A of 40CFR254.521(a)

Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable (a) - Data reported in ug/kg.

DANNELLY ANG - Montgomery, Alabama (*) BACKGROUND INFORMATION 187th Fighter Group

Client Sample ID				B1P-S-0-2	B1P-S-8.5-10	B1P-S-18.5-20	BIP-S-0-2/BIP-S-8.5-10/BIP-S-18.5-20/BIP-S18.5-20D/B3P-S-8.5-10/B3P-S-13.5-15/B3P-S-28.5-30/B5P-S-0-2	B3P-S-8.5-10	B3P-S-13.5-15	B3P-S-28.5-30	B5P-S-0-2
Lab Sample Number				28909001	28909002	28909003	28909004	28922001	28922002	28922003	28909005
Matrix				SOIL	SOIL	SOIL	SOIL	SOIL	TIOS	TIOS	SOIL
Sample Date				2/26/91	33295	2/26/91	2/26/91	2/27/91	16/12/7	16/12/7	2/26/91
ができるとうない。 では、これでは、これでは、これでは、これでは、これでは、これでは、これでは、これ	MCL	CRDL	Action Levels						天工學服務關係關係	A Company of the Assessment of the Company of the C	
Inoganics	ug/L	mg/kg	mg/kg/ug/L					建设制度的	《一》等即译明学	经不完全的 医甲状腺素	2.55年被蒙古
Antimony	9	16.2	30/10	6.2 B	3.9 B	4.0 B		2.7 B	6.1 B	3.3 B	4.0 B
Arsenic	SS	2.7	80/50	11.2	7.1	4.5		5.4	10	7.6	4.8
Barium	2,000	54	4000/1000	51.5 B	39.8 B	29.8 B		239	116	31.6 B	23.2 B
Cadmium	S	1.4	40/10	0.50 B	0.35 B	0.43 B		0.58 B	0.56 B	0.54 B	0.32 B
Chromium	100	2.7	400/50	54.8	15.2	15.2		13.7	20.4	24.9	17.9
Copper	1,300	6.8	400/200	13.3	15.4	18.7		11.6	14.0	23.6	13.7
Lead	15	8.0	N/A/50	37.1	5.4	4.8		6.6	6.7	8.0	8.8
Mercury	2	0.05	20/2.0								
Nickel	100	10.8	2000/100	6.2 B	15.3	15.8		12.8	19.4	20.0	1.9 B
Selenium	20	1.4	N/A/10	0.23 B		0.65 B			0.22 B	0.54 B	
Silver	N/A	2.7	200/50								
Thallium	2	2.7	6/3								0.23 B
Zinc	N/A	5.4	4000/2000	45.0	43.2	46.0		40.0	64.4	62.5	16.6

B - Value detected is less than the CRDL but greater than or equal to the MDL. MCL - Maximum Contaminant Level (U.S. Drinking Water Standards)

CRDL - Contract Required Detection Limit

Biank Space - Value is below the Method Detection Limit(MDL).

CRDL expressed in mg/kg={CRDL in ug/L * Digestion Factor(Assuming 200)
 * Unit Conversion(1/1000)}/%Solids(Assuming 75%)

N/A - Not Applicable
Action Levels proposed in Appendix A of 40CFR254.521(a)
(b) - Soluble Metals
(a) - Data reported in mg/kg (soil) and ug/L (water).

DANNELLY ANG - Montgomery, Alabama (*)

Client Sample ID			-		BSP-S-8.5-10	BSP-S-23.0-25	BSP-S-23.0-25 BSP-S-23.0-25D	BG-MW-6	BG-2P	BG-2P
Lab Sample Number	ber	WWW			28909006	28909007	28909008	18301002	18311002	18311003
Matrix					SOIL	SOIL	SOIL	WATER	WATER	WATER (b)
Sample Date					2/26/91	2/26/91	2/26/91	16/10/4	4/12/91	4/12/91
では、日本の日本の日本の日本の日本の日本の日本の日本の日本の日本の日本の日本の日本の日	MCL	CRDL	CRDL	Action Levels						
Inoganics	ng/L	mg/kg	ug/L	mg/kg/ug/L						
Antimony	9	16.2	09	30/10	5.3 B	4.1 B	4.9 B			
Arsenic	50	2.7	10	80/20	7.5	8.5	10.7			
Barium	2,000	54	200	4000/1000	123	44.4 B	42.6 B			
Cadmium	5	1.4	5	40/10	0.78 B	0.51 B	0.40 B			
Chromium	100	2.7	10	400/50	36.6	24.2	14.4			
Copper	1,300	6.8	25	400/200	13.7	16.4	17.4			
Lead	15	8.0	3	N/A/50	14.2	8.6	11.1			
Mercury	2	0.05	0.2	20/2.0						
Nickel	100	10.8	04	2000/700	20.4	10.1 B	16.2			
Selenium	20	1.4	S	N/A/10		0.58 B	0.49 B			
Silver	N/A	2.7	10	200/50	0.65 B	0.65 B	0.65 B			
Thallium	2	2.7	01	6/3						
Zinc	N/A	5.4	20	4000/2000	36.2	34.6	53.0			
		**************************************	***************************************							

B - Value detected is less than the CRDL but greater than or equal to the MDL. MCL - Maximum Contaminant Level (U.S. Drinking Water Standards)

CRDL - Contract Required Detection Limit

Blank Space - Value is below the Method Detection Limit(MDL).

CRDL expressed in mg/kg={CRDL in ug/L * Digestion Factor(Assuming 200) * Unit Conversion(1/1000)}/%Solids(Assuming 75%)

N/A - Not Applicable
(b) - Soluble Metals
(a) - Data reported in mg/kg (soil) and ug/L (water).

DANNELLY ANG - Montgomery, Alabama " 187th Fighter Group SITE 1, POL AREA

			b c sala	D1BC-8-10	PIRC.8.10D	PIRS.12.14	P2R-S-2-4	P2B-S-8-10	P2B-S-8-10D	P2B-S-12-14	P3B-S-2-4
Client Sample ID			r 153-2-4	01-0-011	201-0-011	2001		2007.000	2007.000	2007 4000	20074000
I ob Somple Number			28998001	28998002	28998003	28998004	28934004	28934005	28934000	78934007	20934000
Metric			SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL
Same Date			3/5/01	3/5/91	3/5/91	3/5/91	2/28/91	2/28/91	2/28/91 2/28/91	2/28/91 2/	2/28/91
Sample Date			2/10/10			en de la constant de	Color and Color and Color of the Letters.	このなるのでは、一日の一日の一日の日本の日本の日本の日本の日本の日本の日本の日本の日本の日本の日本の日本の日本	コテ からに 2000 TME あるののの	とのとなることのできるとのでは、大変の	一一一一一一一一一一一一一一一一一一一一一一一一一一一一一一一一一一一一一一
	CRDL	Action	· · · · · · · · · · · · · · · · · · ·						のでは、大学の大学の大学の大学の大学の大学の大学の大学の大学の大学の大学の大学の大学の大		えることをいうできる。
			20 000 0000000000000000000000000000000	「日本の日本の日本の日本の日本の日本の日本の日本の日本の日本の日本の日本の日本の日	· 中国 图10 10 10 10 10 10 10 10 10 10 10 10 10 1	三年 人名阿里克 医多种 医多种 医多种 医多种	· 多年 · 日本 · 日	古書を見るとは、 は本のは、 は、 は、 は、 は、 は、 は、 は、 は、 は、	方は 法ではおけると	では、一般のないというできた	
BTEX Compounds	ng/kg	ug/kg ug/L							The state of the s		
Benzene	5	N/A/N/A									
Fihylbenzene	5	8E+06/4000									
Toluene	5	2E+07/10,000									
Xvlenes(total)	5	2E+08/70,000		7.9 J							
()	***************************************										

Notes:

J - Estimated (Exceeded holding time)

MCL - Maximum Contaminant Level (U.S. Drinking Water Standards)

CRDL - Contract Required Detection Limit
Action levels proposed in Appendix A of 40CFR253.521(a)
Blank Space - Value is below the Method Detection Limit(MDL).
N/A - Not Applicable
(a) - Data reported in ug/kg (soil) and ug/L (water).

187th Fighter Group SITE 1, POL AREA

DANNELLY ANG - Montgomery, Alabama (*)

Client Sample ID					P3B-S-6-8	P3B-S-12-14	P4BS-2-4	P4BS-6-8	P4BS-12-14	P5B-S-4-6	P3B-S-6-8 P3B-S-12-14 P4BS-2-4 P4BS-6-8 P4BS-12-14 P5B-S-4-6 P5B-S-12-14 P5B-S-8-10 P5BW1	P5B-S-8-10	P5BW1
Lab Sample Number					28934009	28934010 28998011 28998012 28934011	28998011	28998012	28998013	28934011	28934013	28934012 2897200	28972001
Matrix					SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	Г	SOIL	WATER
Sample Date					2/28/91	2/28/91	3/5/91	3/2/61	3/5/91	2/28/91	2/28/91	2/28/91	3/4/91
	MCL	MCL CRDL CRDL	CRDL	Action							できる世界を選集	(A)	は発音される
BTEX Compounds	ug/L	ug/L ug/kg ug/L	ng/L	ug/kg ug/L									
Benzene	5	5	5	N/A/N/A									400
Ethylbenzene	700	5	5	8E+06/4000			12	6.7		3200		340	15
Toluene	1,000	5	5	2E+07/10,000									120
Xylenes(total)	10,000	5	5	2E+08/70,000		6.9	34			3000		1100	260

J - Estimated (Exceeded holding time)

MCL - Maximum Contaminant Level (U.S. Drinking Water Standards)

CRDL - Contract Required Detection Limit

Blank Space - Value is below the Method Detection Limit(MDL). Action levels proposed in Appendix A of 40CFR253.521(a)

N/A - Not Applicable (a) - Data reported in ug/kg (soil) and ug/L (water).

DANNELLY ANG - Montgomery, Alabama " 187th Fighter Group SITE 1, POL AREA

								0 / 0454	01 0 3020	11 61 9000	Wara	D8RC-2-4	P8BS-6-8	P8BS-12-14
Cliont Sumalo II)					8-9-SE94	DeBS-8-10	P6BS-12-14	P/BS-0-8	F/BS-6-10	F/D3-12-14	11011	7.000		7100000
Cilcin Sample 10	1				30000000	20090000	70080080	28008008	28998009	28998010	29041002	28998014	28998015	28998016
1 sh Sample Number					C0006607	2022000	1000/107	20000				100	1100	шов
					TO9	поз	SOIL	SOIL	SOIL	SOIL	WAIEK	SOIL	SOIL	3010
Matrix					3000				101310	275/01	377/91	1/5/01	16/5/2	3/5/91
-					3/2/01	3/2/9]	3/5/91	3/3/91	16/6/6	16/0/6	3/1/21	171010		
Sample Date	_	1				Company of the Compan		Control of the Contro	の一の日本の大学の日本の日本の日本の日本の日本の日本の日本の日本の日本の日本の日本の日本の日本の	PARTY OF SERVICE AND ADDRESS OF THE PARTY OF	二、大学には はない はない はない はない はない はない はない はない はない はな	このであるとはあるというである	とを見れるというないとなっている	今年 できたの言意を向け
	I VV	Juan	ותמט	Action								The state of the s		
The second secon	1	1						SHIP CHANGE BY SELECTION OF THE SECURIOR SHAPE S	· · · · · · · · · · · · · · · · · · ·	CONTRACTOR STREET	がからの製造の日本で、名前の	代記は各体を持ちたです。	これでは あるがに	おないないない を記
DTCV Compounds	110/1	110/kg	ne/L	ne/ke ue/L				THE PROPERTY OF STREET			SECTION REPRESENTATION OF STREET	desired to the collective properties as which the last		
Dica Compounds		4 4												
Renzene	٠	'n	'n	Z/Z/Z/Z										
a constant	3		,	00//VV					4.0					
Ethylbenzene	3	2		SE-FOOT-FOOD										
	8	_	~	2F±07/10 000	_	_								
loidene	1,000	,		20101				3,			03			_
Videnac/total)	000	v	S	2E+08/70.000				13			7	,		
A VICINES (10tm)	22.5													

Notes:

J. Estimated (Exceeded holding time)

MCL. Maximum Contaminant Level (U.S. Drinking Water Standards)

MCL. Contract Required Detection Limit

Action levels proposed in Appendix A of 40CFR253.521(a)

Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

(a) - Data reported in ug/kg (soil) and ug/L (water).

2/26/91 **B5PS** 17937 SOIL 10 DANNELLY ANG - Montgomery, Alabama (a) 2/26/91 B3PS 17931 SOIL 28.4 2/26/91 B1PS 17937 SOIL 4.8 3.9 3.6 **Fotal Petroleum Hydrocarbons** Sample Depth(Feet) 0-2 ab Sample Number 13.5-15 18.5-20 23-25 28.5-30 8.5-10 Client Sample ID Sample Date Matrix

Notes:

Shaded areas indicate depth intervals not sampled.

DANNELLY ANG - Montgomery, Alabama (a) 187th Fighter Group SITE 1, POL AREA

Lab Sample Number Matrix Matrix Sample Date Sample Date CRDL Action PAH Compounds ug/kg Levels 1-Methylnaphthalene 330 N/A 2-Methylnaphthalene 330 N/A		24 2 224	701-0-0711 01-0-0711	+1-71-CG11	L7D-2-7-4		P2B-S-8-10 P2B-S-8-10D
CRDL ug/kg 330 330	28998001	28998002	28998003	28998004	28934004	28934005	28934006
CRDL ug/kg 330 330	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL
CRDL ug/kg 330 330	3/5/91	3/2/61	3/5/91	3/5/91	2/28/91	2/28/91	2/28/91
1330 330	u(
330						Assessment of the second	Target and Property Park
330	1						
Naphthalene 330 N/A	1						
Acenaphthylene 330 N/A	1						
Acenaphthene 330 N/A	1						
Fluorene 330 N/A							
Phenanthrene 330 N/A							
Fluoranthene 330 N/A	1						
Pyrene 330 N/A							

J - Estimated (Exceeded holding time)

CRDL - Contract Required Detection Limit

Blank Space - Value is below the Method Detection Limit(MDL). Action Levels are proposed in Appendix A of 40CFR252.521(a)

N/A - Not Applicable

(a) - Data reported in ug/kg (soil) and ug/L (water).

DANNELLY ANG - Montgomery, Alabama (*) 187th Fighter Group SITE 1, POL AREA

Client Somule ID			P1BS-2-4	P1BS-8-10	P1BS-8-10D	P1BS-12-14	P2B-S-2-4	P2B-S-8-10	P2B-S-8-10D
Lab Sample Number			28998001	28998002	28998003	28998004	28934004	28934005	28934006
Matrix			SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL
Sample Date			3/5/91	3/5/91	3/5/91	3/5/91	2/28/91	2/28/91	2/28/91
1000年の日本の日本の日本の日本の日本の日本の日本の日本の日本の日本の日本の日本の日本の	CRDL	Action							なる。
PAH Compounds	ug/kg	Levels							
Benzo(a)anthracene	330	N/A		·					
Chrysene	330	N/A							
Benzo(b)fluoranthene	330	N/A							
Benzo(k)fluoranthene	330	N/A							
Benzo(a)pyrene	330	N/A							
Indeno(1,2,3-cd)pyrene	330	N/A							
Dibenzo(a,h)anthracene	330	N/A							
Benzo(g,h,i)perylene	330	N/A							

J - Estimated (Exceeded holding time)

CRDL - Contract Required Detection Limit Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

Action Levels are proposed in Appendix A of 40CFR254.521(a) (a) - Data reported in ug/kg (soil) and ug/L (water).

DANNELLY ANG - Montgomery, Alabama (*) 187th Fighter Group SITE 1, POL AREA

Client Sample ID			P2B-S-12-14	P3B-S-2-4	P3B-S-6-8	P3B-S-12-14	P4BS-2-4	P4BS-6-8	P4BS-12-14
Lab Sample Number			28934007	ı	28934009	28934010	28998011	28998012	28998013
Matrix			SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL
Sample Date			2/28/91	2/28/91	2/28/91	2/28/91	3/5/91	3/5/91	3/5/91
	CRDL	Action		· · · · · · · · · · · · · · · · · · ·		· · · · · · · · · · · · · · · · · · ·	1. 不是不管情况下了		
PAH Compounds	ug/kg	Levels					Ander Williams Wer		
1-Methylnaphthalene	330	N/A							
2-Methylnaphthalene	330	N/A							
Naphthalene	330	N/A							
Acenaphthylene	330	N/A							
Acenaphthene	330	N/A							
Fluorene	330	N/A							
Phenanthrene	330	N/A							
Fluoranthene	330	N/A							
Pyrene	330	N/A							

J - Estimated (Exceeded holding time)

CRDL - Contract Required Detection Limit Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

Action Levels are proposed in Appendix A of 40CFR254.521(a) (a) - Data reported in ug/kg (soil) and ug/L (water).

187th Fighter Group SITE 1, POL AREA

DANNELLY ANG - Montgomery, Alabama (*)

Client Sample ID			P1BS-2-4	P1BS-8-10	P1BS-8-10D	P1BS-12-14	P2B-S-2-4	P2B-S-8-10	P2B-S-8-10D
Lab Sample Number			28998001	28998002	28998003	28998004	28934004	28934005	28934006
Matrix			SOIL	SOIL	SOIL	TIOS	SOIL	SOIL	SOIL
Sample Date			3/5/91	16/5/8	3/5/91	3/5/91	2/28/91	2/28/91	2/28/91
The second secon	CRDL	Action						- 一直の一直の	
PAH Compounds	ug/kg	Levels	[1988] 图 图 2018 图 2018	多為 的特殊 網 的 群群				高级的数据等 对	· · · · · · · · · · · · · · · · · · ·
Benzo(a)anthracene	330	N/A							
Chrysene	330	N/A							
Benzo(b)fluoranthene	330	N/A							
Benzo(k)fluoranthene	330	N/A		,					
Benzo(a)pyrene	330	N/A							
Indeno(1,2,3-cd)pyrene	330	N/A							
Dibenzo(a,h)anthracene	330	N/A							
Benzo(g,h,i)perylene	330	N/A							

J - Estimated (Exceeded holding time) CRDL - Contract Required Detection Limit

Blank Space - Value is below the Method Detection Limit(MDL). N/A - Not Applicable Action Levels are proposed in Appendix A of 40CFR254.521(a) (a) - Data reported in ug/kg (soil) and ug/L (water).

DANNELLY ANG - Montgomery, Alabama (a) 187th Fighter Group SITE 1, POL AREA

Oliver Secure 10					P5B-S-4-6	P5B-S-8-10	P5B-S-12-14	P5BW1	P6BS-6-8	P6BS-8-10	P6BS-12-14
r ob Somele Number					28934011	28934012	28934013	28972001	28998005	28998006	28998007
Motrix					SOIL	SOIL	TIOS	WATER	SOIL	SOIL	SOIL
Sample Date					2/28/91	2/28/91	2/28/91	3/4/91	3/5/91	3/5/91	3/5/91
	MCL	CRDL	CRDL	Action							の一般の
PAH Compounds	ng/L	ug/kg	ug/L	Levels				一番 の 日本 の 日	できるのできる数を重要		
1-Methylnaphthalene	A/A	330	10	N/A				200			
2-Methylnaphthalene	N/A	330	10	N/A				250			
Naphthalene	N/A	330	10	N/A				87			
Acenaphthylene	A/A	330	10	N/A							
Acenaphthene	A/A	330	10	N/A							
Fluorene	A/A	330	10	N/A							
Phenanthrene	A/A	330	10	N/A							
Fluoranthene	N/A	330	10	N/A							
Pyrene	N.A	330	10	N/A							

Notes:

J - Estimated (Exceeded holding time)

MCL - Maximum Contaminant Level (U.S. Drinking Water Standards)

CRDL - Contract Required Detection Limit

Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

Action Levels are proposed in Appendix A of 40CFR254,521(a)

(a) - Data reported in ug/kg (soil) and ug/L (water).

DANNELLY ANG - Montgomery, Alabama (a) 187th Fighter Group SITE 1, POL AREA

Client Sample ID					P5B-S-4-6	P5B-S-8-10	P5B-S-12-14	P5BW1	P6BS-6-8	P6BS-8-10	P6BS-12-14
Lab Sample Number					28934011	28934012	28934013	28972001	28998005	28998006	28998007
Matrix					SOIL	SOIL	SOIL	WATER	SOIL	SOIL	SOIL
Sample Date					2/28/91	2/28/91	2/28/91	3/4/91	3/5/91	3/5/91	3/5/91
	MCL	CRDL	CRDL	Action			10 17 18 18 14 15 18 18 18 18 18 18 18 18 18 18 18 18 18				
PAH Compounds	ng/L	ug/kg	ng/L	Levels							
Benzo(a)anthracene	0.1	330	10	N/A							
Chrysene	0.2	330	10	N/A							
Benzo(b)fluoranthene	0.2	330	10	N/A							
Benzo(k)fluoranthene	0.7	330	10	N/A							
Benzo(a)pyrene	0.7	330	10	N/A							
Indeno(1,2,3-cd)pyrene	0.4	330	10	N/A							
Dibenzo(a,h)anthracene	0.3	330	10	N/A							
Benzo(g,h,i)perylene	N/A	330	10	N/A							

J - Estimated (Exceeded holding time)
MCL - Maximum Contaminant Level (U.S. Drinking Water Standards)

CRDL - Contract Required Detection Limit
Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable
Action Levels are proposed in Appendix A of 40CFR254.521(a)

(a) - Data reported in ug/kg (soil) and ug/L (water).

DANNELLY ANG - Montgomery, Alabama (*) 187th Fighter Group SITE 1, POL AREA

Client Sample ID					P7BS-6-8	P7BS-8-10	P7BS-12-14	P7BW	P8BS-2-4	P8BS-6-8	P8BS-12-14
Lab Sample Number					28998008	28998009	28998010	29041002	28998014	28998015	28998016
Matrix					SOIL	SOIL	SOIL	WATER	SOIL	SOLL	SOIL
Sample Date					3/5/91	3/5/91	3/5/91	3/7/91	3/5/91	3/5/91	3/5/91
「おける」というでは、またのでは、これは、おけるのであった。	MCL	CRDL	CRDL	Action							日本のでは、日本の
PAH Compounds	ng/L	ug/kg	ug/L	Levels							
1-Methylnaphthalene	N/A	330	10	N/A				130			
2-Methylnaphthalene	N/A	330	10	N/A				7			
Naphthalene	A/N	330	10	N/A							
Acenaphthylene	N/A	330	10	N/A							
Acenaphthene	N/A	330	10	N/A							
Fluorene	N/A	330	10	N/A							
Phenanthrene	N/A	330	10	N/A							
Fluoranthene	N/A	330	10	N/A							
Pyrene	N.A	330	10	N/A							

J - Estimated (Exceeded holding time)
MCL - Maximum Contaminant Level (U.S. Drinking Water Standards)
CRDL - Contract Required Detection Limit
Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable
Action Levels are proposed in Appendix A of 40CFR254.521(a)
(a) - Data reported in ug/kg (soil) and ug/L (water).

DANNELLY ANG - Montgomery, Alabama (*) 187th Fighter Group SITE 1, POL AREA

Client Sample ID					P7BS-6-8	P7BS-8-10	P7BS-12-14	P7BW	P8BS-2-4	P8BS-6-8	P8BS-12-14
Lab Sample Number					28998008	28998009	28998010	29041002	28998014	28998015	28998016
Matrix					SOIL	SOIL	SOIL	WATER	SOIL	SOIL	SOIL
Sample Date					3/5/91	3/5/91	3/5/91	3/7/91	3/5/91	3/5/91	3/5/91
三種の天然を発がれた。 ならの 古屋のからと	MCL	CRDL	CRDL	Action						和海洲山海河和北京河南	
PAH Compounds	ug/L	ug/kg	ug/L	Levels					· · · · · · · · · · · · · · · · · · ·	東京の日本の日本の日本の日本の日本の日本の日本の日本の日本の日本の日本の日本の日本の	三、大学の大学の大学で
Benzo(a)anthracene	0.1	330	10	N/A							
Chrysene	0.2	330	10	N/A							
Benzo(b)fluoranthene	0.2	330	10	N/A					·		
Benzo(k)fluoranthene	0.2	330	10	N/A							
Benzo(a)pyrene	0.7	330	10	N/A							
Indeno(1,2,3-cd)pyrene	0.4	330	10	N/A							
Dibenzo(a,h)anthracene	0.3	330	10	N/A							
Benzo(g,h,i)perylene	N/A	330	10	N/A							

J - Estimated (Exceeded holding time)
MCL - Maximum Contaminant Level (U.S. Drinking Water Standards)
CRDL - Contract Required Detection Limit
Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

Action Levels are proposed in Appendix A of 40CFR254.521(a) (a) - Data reported in ug/kg (soil) and ug/L (water).

DANNELLY ANG - Montgomery, Alabama (*) SITE 2, OIL WATER SEPARATOR 187th Fighter Group

Client Somule ID			03BS-12-14	03BS-12-14D	Ł	03BS12-14 MS 03BS12-14 DUP	04BS-0-2	04BS-6-8	04BS-14-16	04BS-14-16 04BS-14-16D
I ah Sample Number			29017013	29017014		29017P14	29017001	29017002	29017003	29017004
Matrix			SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL
Sample Date			3/6/91	3/6/91	3/6/91	3/6/91	3/6/91	3/6/91	3/6/91	3/6/91
を記した。なるとは、世上の人が記していた。 では、一般には、一般には、一般には、一般には、一般には、一般には、一般には、一般に	CRDL	Action								
Volatile Compounds	ug/kg	Levels					ですが、おきを選挙を発言	このでは、 一般のない		
1.1.2.2-Tetrachloroethane	5	40,000								
1,1-Dichloroethane	5	N/A								790
1.1-Dichloroethene	5	N/A								
1.2-Dichloroethene(total)	5	8,000	730 J	660 J				5400	220 J	096
Acetone	10	8E+06	2900 B	2900 B			5000 BJ	13000 B	4800 B	1800 B
Benzene	5	N/A					1000 J			
Carbon Disulfide	5	8E+06								
Ethylbenzene	5	8E+06		,			11000	1300 J		
Methylene Chloride	5	90,000	3600 B	2900 B			9500 B	7900 B		
Tetrachloroethene	5	N/A							400 J	1100

B - Applies to organic data only. Present in the corresponding method blank
J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.
E - Applies to Gas Chromatography/Mass Spectroscopy data only.
Indicate compound above or below linear range of instrument.

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a)

Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

(a) - Data reported in ug/kg.

SITE 2, OIL WATER SEPARATOR 187th Fighter Group

DANNELLY ANG - Montgomery, Alabama (*)

e Number SOIL 29017013 29017014 29017M14 29017P14 29017001 te SOIL SO	Client Sample ID			03BS-12-14		03BS12-14 MS	03BS-12-14D 03BS12-14 MS 03BS12-14 DUP	04BS-0-2	04BS-6-8	04BS-14-16
CRDL Action SOIL <	Lab Sample Number			29017013	29017014	29017M14	29017P14	29017001	29017002	29017003
CRDL Action 3/6/91 <td>Matrix</td> <td></td> <td></td> <td>SOIL</td> <td>SOIL</td> <td>SOIL</td> <td>SOIL</td> <td>SOIL</td> <td>SOIL</td> <td>SOIL</td>	Matrix			SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL
CRDL Action Evels Evels <th< td=""><td>Sample Date</td><td></td><td></td><td>3/6/91</td><td>3/6/91</td><td>3/6/91</td><td>3/6/91</td><td>3/6/91</td><td>3/6/91</td><td>3/6/91</td></th<>	Sample Date			3/6/91	3/6/91	3/6/91	3/6/91	3/6/91	3/6/91	3/6/91
pounds ug/kg Levels Tevels Levels Levels </td <td></td> <td>CRDL</td> <td>Action</td> <td>计算中部经验和印象对</td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td>		CRDL	Action	计算中部经验和印象对						
ne 5 2E+08 180 J 20000 21000 le 10 N/A 27000 10000 89000 1 le 10 N/A 89000 89000 1 le 10 N/A 89000 1 le 10 N/A 10 10,000 1 lhane 5 100,000 1 1 1 lhane 5 N/A 1 1 1	Volatile Compounds	ug/kg	Levels		医多种类型核酮 酸器	多層等數學學的		· 100 · 100		
ne 5 N/A 27000 10000 6 89000 ie 10 N/A 89000 89000 89000 ic 10 N/A 89000 89000 89000 89000 ic 10 N/A 890000 89000 89000 89000	Toluene	5	2E+08	180 J				21000	5300	310 J
le 10 N/A 89000 se 10 N/A 89000 sthane 5 N/A 89000 shape 89000 89000 89000	Trichloroethene	5	N/A	27000	10000					33000
se 2E+08 89000 ne 10 N/A 89000 ne 10 100,000 10 s 100,000 100,000 100,000 ithane 5 N/A 10	Vinyl Chloride	10	N/A							
ane 10 ane 10 le 10 cethane 5	Xylene(total)	5	2E+08					89000	17000	
ane 10 le 10 sorthane 5	Chloromethane	10	N/A							
te 10 5 5 5 5 10 10 10 10 10 10 10 10 10 10 10 10 10	Bromomethane	10	100,000							
5 Dethane 5	Chloroethane	10	N/A							
oethane 5	Chloroform	5	100,000							
10	1,2-Dichloroethane	5	N/A							
	2-Butanone	10	N/A							

Notes:

- B Applies to organic data only. Present in the corresponding method blank
- J Applies to organic data only. Value detected is greater than zero but less than the CRDL.
 - E Applies to Gas Chromatography/Mass Spectroscopy data only.

Indicate compound above or below linear range of instrument.

Action Levels proposed in Appendix A of 40CFR254.521(a) CRDL - Contract Required Detection Limit

Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

DANNELLY ANG - Montgomery, Alabama 🗝 SITE 2, OIL WATER SEPARATOR 187th Fighter Group

Client Sample ID			03BS-12-14	03BS-12-14D	03BS12-14 MS	03BS-12-14 03BS-12-14D 03BS12-14 MS 03BS12-14 DUP	04BS-0-2	04BS-6-8	04BS-14-16
Lab Sample Number			29017013	29017014	29017M14	29017P14	29017001	29017002	29017003
Matrix			SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL
Sample Date			3/6/91	3/6/91	3/6/91	3/6/91	3/6/91	3/6/91	3/6/91
	CRDL	Action							
Volatile Compounds	ug/kg	Levels							74.5 24.9 24.9 2.0 2.0 2.0 2.0 2.0 2.0 2.0 2.0 2.0 2.0
1,1,1-Trichloroethane	5	7E+06							
Carbon Tetrachloride	5	1,000							
Vinyl Acetate	10	N/A							
Bromodichloromethane	5	200							
1,2-Dichloropropane	5	N/A							
cis-1,3-Dichloropropene	5	20,000							
Dibromochloromethane	5	N/A							
1,1,2-Trichloroethane	5	100,000							
trans-1,3-Dichloropropene	5	20,000							
Bromoform	5	2E+06							

- B Applies to organic data only. Present in the corresponding method blank
 J Applies to organic data only. Value detected is greater than zero but less than the CRDL.
 E Applies to Gas Chromatography/Mass Spectroscopy data only.
 Indicate compound above or below linear range of instrument.

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a) Blank Space - Value is below the Method Detection Limit(MDL). N/A - Not Applicable

DANNELLY ANG - Montgomery, Alabama (*)

Client Sample ID			03BS-12-14	03BS-12-14D	03BS12-14 MS	03BS-12-14D 03BS12-14 MS 03BS12-14 DUP	04BS-0-2	04BS-6-8	04BS-14-16
Lab Sample Number			29017013	29017014	29017M14	29017P14		29017002	29017003
Matrix			SOIL	TIOS	SOIL	SOIL	SOIL	SOIL	SOIL
Sample Date			3/6/91	3/6/91	3/6/91	3/6/91	3/6/91	3/6/91	3/6/91
	CRDL	Action							
Volatile Compounds	ug/kg	Levels							
4-Methyl-2-Pentanone	10	N/A							
2-Hexanone	10	N/A							
Chlorobenzene	5	2E+06							
Styrene	5	2E+07							

Notes:

B - Applies to organic data only. Present in the corresponding method blank J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.

E - Applies to Gas Chromatography/Mass Spectroscopy data only.

Indicate compound above or below linear range of instrument.

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a)

Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

DANNELLY ANG - Montgomery, Alabama (*) SITE 2, OIL WATER SEPARATOR 187th Fighter Group

			חזו זו שמזט	600000	0500 0 10	05BC 12 14
Cilent Sample ID			04D2-14-10D	7-0-00C0	07-0-000	17-17-0000
Lab Sample Number		;	29017004	29017008	29017009	29017010
Matrix			SOIL	SOIL	SOIL	SOIL
Sample Date			3/6/91	3/6/91	3/6/91	3/6/91
	CRDL	Action				
Volatile Compounds	ug/kg	Levels				
1,1,2,2-Tetrachloroethane	. 5	40,000				
1,1-Dichloroethane	5	N/A	190			
1,1-Dichloroethene	5	N/A				
1,2-Dichloroethene(total)	5	8,000	096		28000	
Acetone	10	8E+06	1800 B	2900	2000	
Benzene	5	N/A		390 J		
Carbon Disulfide	5	90+ 3 8				
Ethylbenzene	5	90+ 3 8		2300	220 J	
Methylene Chloride	5	000'06		3700 B	810 J	2600 B
Tetrachloroethene	5	N/A	1100		·	

Notes:

- B Applies to organic data only. Present in the corresponding method blank
- J Applies to organic data only. Value detected is greater than zero but less than the CRDL.
 - E Applies to Gas Chromatography/Mass Spectroscopy data only.

Indicate compound above or below linear range of instrument.

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a) Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

DANNELLY ANG - Montgomery, Alabama (a)

Client Sample ID			04BS-14-16D	05BS-0-2	05BS-8-10	05BS-12-14
Lab Sample Number			29017004	29017008	29017009	29017010
Matrix			TIOS	SOIL	TIOS	SOIL
Sample Date			3/6/91	3/6/91	3/6/91	3/6/91
	CRDL	Action				
Volatile Compounds	ug/kg	Levels	1.46.1的14.50.144.744.744			
Toluene	5	2E+08	1100	5500	640 J	
Trichloroethene	5	N/A	65000 E		4000	13000
Vinyl Chloride	10	N/A			1300 J	
Xylene(total)	. 5	2E+08	790 J	17000	1200	
Chloromethane	10	N/A				
Bromomethane	10	100,000				
Chloroethane	10	N/A				
Chloroform	5	100,000				
1,2-Dichloroethane	5	N/A		,		
2-Butanone	10	N/A				

Notes:

B - Applies to organic data only. Present in the corresponding method blank

J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.

E - Applies to Gas Chromatography/Mass Spectroscopy data only.

Indicate compound above or below linear range of instrument.

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a)

Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

(a) - Data reported in ug/kg.

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DANNELLY ANG - Montgomery, Alabama (a)

Client Sample ID			04BS-14-16D	05BS-0-2	05BS-8-10	05BS-12-14
Lab Sample Number			29017004		29017009	29017010
Matrix			SOIL	SOIL	SOIL	SOIL
Sample Date			3/6/91	3/6/91	3/6/91	3/6/91
	CRDL	Action				
Volatile Compounds	ug/kg	Levels		7		
1,1,1-Trichloroethane	5	7E+06				
Carbon Tetrachloride	5	1,000				
Vinyl Acetate	10	N/A				
Bromodichloromethane	5	200				
1,2-Dichloropropane	5	N/A			•	
cis-1,3-Dichloropropene	5	20,000				
Dibromochloromethane	5	N/A				
1,1,2-Trichloroethane	5	100,000				
trans-1,3-Dichloropropene	5	20,000				
Bromoform	5	2E+06				

Notes.

- B Applies to organic data only. Present in the corresponding method blank
- J Applies to organic data only. Value detected is greater than zero but less than the CRDL.
 - E Applies to Gas Chromatography/Mass Spectroscopy data only. Indicate compound above or below linear range of instrument.

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a)

Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

DANNELLY ANG - Montgomery, Alabama (a)

Client Sample ID			04BS-14-16D	05BS-0-2	05BS-8-10	
Lab Sample Number		-	29017004	29017008	29017009	29017010
Matrix			SOIL	SOIL	SOIL	
Sample Date			3/6/91	3/6/91	3/6/91	3/6/91
	CRDL	Action				
Volatile Compounds	ug/kg	Levels				张文明学科 中国学科
4-Methyl-2-Pentanone	10	N/A				
2-Hexanone	10	N/A				
Chlorobenzene	5	2E+06				
Styrene	5	2E+07				

Notes:

B - Applies to organic data only. Present in the corresponding method blank

J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.

E - Applies to Gas Chromatography/Mass Spectroscopy data only.

Indicate compound above or below linear range of instrument.

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a)

Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

SITE 2, OIL WATER SEPARATOR 187th Fighter Group DANNELLY ANG - Montgomery, Alabama (*)

Client Sample ID			01BS-0-2	01BS-10-12	01BS-14-16	02B-S-0-2	01BS-14-16 02B-S-0-2 02B-S-12-14 02B-S-18-20		03BS-0-2	03BS-6-8
Lab Sample Number			29017005	29017006	29017007	28934001	28934002	28934003	29017011	29017012
Matrix			SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL
Sample Date			3/6/91	3/6/91	3/6/91	2/28/91	2/28/91	2/28/91	3/6/91	3/6/91
	CRDL	Action								
Volatile Compounds	ug/kg	Levels							STREET, WHILE THE	
1,1,2,2-Tetrachloroethane	5	40,000			1600			,		
1,1-Dichloroethane	5	N/A		1 J						
1,1-Dichloroethene	5	N/A						24 J		
1,2-Dichloroethene(total)	5	8,000		180	1500		150	1000	700	830
Acetone	10	8E+06	53	19 B	1900 B		17 B	100 B	30 BJ	1900
Benzene	5	N/A	3 J							
Carbon Disulfide	5	8E+06						18 J		
Ethylbenzene	5	8E+06	15							
Methylene Chloride	5	000'06	14 B	11 B	2000	13 B	21 B	120 B	51 B	1300 J
Tetrachloroethene	5	N/A			450 J					

Notes

B - Applies to organic data only. Present in the corresponding method blank

J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.

Applies to Gas Chromatography/Mass Spectroscopy data only.
 Indicate compound above or below linear range of instrument.

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a)

Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

DANNELLY ANG - Montgomery, Alabama (*)

Client Sample ID			01BS-0-2	01BS-10-12	01BS-14-16	02B-S-0-2	01BS-10-12 01BS-14-16 02B-S-0-2 02B-S-12-14 02B-S-18-20	02B-S-18-20	03BS-0-2	03BS-6-8
Lab Sample Number			29017005	29017006	29017007	28934001	28934002	28934003	29017011	29017012
Matrix			SOIL	SOIL	SOIL	SOIL	SOIL	TIOS	SOIL	SOIL
Sample Date			3/6/91	3/6/91	3/6/91	2/28/91	2/28/91	2/28/91	3/6/91	3/6/91
	CRDL	Action	训练表现。 "哈里"		心影声级影响					
Volatile Compounds	ug/kg	Levels			有多一个商务	30000000000000000000000000000000000000	· 第2、2000年 12、5年			
1,1,1-Trichloroethane	5	7E+06								
Carbon Tetrachloride	5	1,000								
Vinyl Acetate	10	N/A								
Bromodichloromethane	5	200								
1,2-Dichloropropane	5	N/A								
cis-1,3-Dichloropropene	5	20,000								
Dibromochloromethane	5	N/A								
1,1,2-Trichloroethane	5	100,000								
trans-1,3-Dichloropropene	5	20,000								
Bromoform	5	2E+06								

B - Applies to organic data only. Present in the corresponding method blank
 J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.
 E - Applies to Gas Chromatography/Mass Spectroscopy data only.

Indicate compound above or below linear range of instrument. CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a)

Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

DANNELLY ANG - Montgomery, Alabama (a)

Client Somple ID			01BS-0-2	01BS-10-12	01BS-14-16	02B-S-0-2	02B-S-12-14	01BS-10-12 01BS-14-16 02B-S-0-2 02B-S-12-14 02B-S-18-20	03BS-0-2	03BS-6-8
I ob Sample Number			29017005	29017006	29017007	28934001	28934002	28934003	29017011	29017012
Matrix			SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL
Sample Date			3/6/91	3/6/91	3/6/91	2/28/91	2/28/91	2/28/91	3/6/91	3/6/91
がは、これのでは、一般のでは、これの	CRDL	Action								
Volatile Compounds	ug/kg	Levels		· · · · · · · · · · · · · · · · · · ·					然於美麗樓等	
4-Methyl-2-Pentanone	10	N/A								
2-Hexanone	10	N/A								
Chlorobenzene	5	2E+06								
Styrene	5	2E+07								

Notes:

B - Applies to organic data only. Present in the corresponding method blank

J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.

E - Applies to Gas Chromatography/Mass Spectroscopy data only.

Indicate compound above or below linear range of instrument.

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a)

Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

DANNELLY ANG - Montgomery, Alabama (*)

Client Sample ID			01BS-0-2	01BS-10-12	01BS-14-16	02B-S-0-2	01BS-10-12 01BS-14-16 02B-S-0-2 02B-S-12-14 02B-S-18-20 03BS-0-2	02B-S-18-20	03BS-0-2	03BS-6-8
Lab Sample Number			29017005	29017006	29017007	28934001	28934002	28934003	29017011	29017012
Matrix			SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL
Sample Date			3/6/91	3/6/91	3/6/91	2/28/91	2/28/91	2/28/91	3/6/91	3/6/91
	CRDL	Action								1. SAN TO 12
Semivolatile Compounds	ug/kg	Levels (ug/kg)								3. 情况数 (新水石)
2-Methylnaphthalene	330	N/A								
Anthracene	330	N/A					120 J			
Benzo(a)anthracene	330	N/A				f 69	170			
Benzo(a)pyrene	330	N/A		,		75 J	099			
Benzo(b)fluoranthene	330	N/A				96 J	069			
Benzo(g,h,i)perylene	330	N/A				60 J	440			
Benzo(k)fluoranthene	330	N/A				94 J	092			
Benzoic Acid	1600	N/A								
Chrysene	330	N/A				94 J	780			
Di-n-butylphthalate	330	8E+06				42 J				

Notes:

B - Applies to organic data only. Present in the corresponding method blank

J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.

E - Applies to Gas Chromatography/Mass Spectroscopy data only.

Indicate compound above or below linear range of instrument. CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a)

Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

DANNELLY ANG - Montgomery, Alabama (*)

			,					00 01 0 200	0 0 0 0 0 0 0	0 / 0000
Client Sample ID	-		01BS-0-2	01BS-10-12	01BS-10-12 01BS-14-16	02B-S-0-2	02B-S-12-14	02B-S-18-20	03BS-0-2	U3B3-0-8
Lab Sample Number			29017005	29017006	29017007	28934001	28934002	28934003	29017011	29017012
Matrix			SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL
Sample Date			3/6/91	3/6/91	3/6/91	2/28/91	2/28/91	2/28/91	3/6/91	3/6/91
· · · · · · · · · · · · · · · · · · ·	CRDL	Action	· · · · · · · · · · · · · · · · · · ·							一年 ののはないのではない
Semivolatile Compounds	ug/kg	Levels (ug/kg)						である。		公司等 化新国家特别的
Dibenz(a,h)anthracene	330	N/A					120 J			
Fluoranthene	330	N/A	•			140 J	1600			
Fluorene	330	N/A								
Indeno(1,2,3-cd)pyrene	330	N/A				56 J	400 J			
N-Nitrosodiphenylamine	330	100,000					43 J			
Naphthalene	330	N/A								
Nitrobenzene	330	40,000								
Phenanthrene	330	N/A					510			
Pyrene	330	N/A				110 J	1200			
Bis(2-Ethylhexyl)phthalate	330	50,000	310 BJ	2900 B	1000 B	3000 B	2800 B	1900 B	440 B	3900 B

Note

- B Applies to organic data only. Present in the corresponding method blank
- J Applies to organic data only. Value detected is greater than zero but less than the CRDL.
 - E Applies to Gas Chromatography/Mass Spectroscopy data only.
 - Indicate compound above or below linear range of instrument.
 - CRDL Contract Required Detection Limit
- Action Levels proposed in Appendix A of 40CFR254.521(a)
- Blank Space Value is below the Method Detection Limit(MDL).
- N/A Not Applicable
- (a) Data reported in ug/kg.

DANNELLY ANG - Montgomery, Alabama (*)

Client Sample ID			01BS-0-2	01BS-10-12	01BS-14-16	02B-S-0-2	02B-S-12-14	01BS-10-12 01BS-14-16 02B-S-0-2 02B-S-12-14 02B-S-18-20	03BS-0-2	03BS-6-8
Lab Sample Number			29017005	29017006	29017007	28934001	28934002	28934003	29017011	29017012
Matrix			SOIL	SOIL	SOIL	SOIL	TIOS	SOIL	SOIL	SOIL
Sample Date			3/6/91	3/6/91	3/6/91	2/28/91	2/28/91	2/28/91	3/6/91	3/6/91
	CRDL	Action	。 			以外,即将该外				
Semivolatile Compounds	ug/kg	Levels (ug/kg)		開発開開機器系統第		新物学用学用的				
Phenol	330	5E+07								
bis(2-Chloroethy1)Ether	330	N/A								-
2-Chlorophenol	330	400,000								
1,3-Dichlorobenzene	330	N/A								
1,4-Dichlorobenzene	330	N/A								
Benzyl Alcohol	330	N/A								
1,2-Dichlorobenzene	330	N/A								
2-Methylphenol	330	N/A								
bis(2-Chloroisopropyl)Ether	330	N/A								
4-Methylphenol	330	N/A								

Notes:

- B Applies to organic data only. Present in the corresponding method blank
- J Applies to organic data only. Value detected is greater than zero but less than the CRDL.
 - E Applies to Gas Chromatography/Mass Spectroscopy data only.

Indicate compound above or below linear range of instrument.

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a) Blank Space - Value is below the Method Detection Limit(MDL).

Blank Space - Value is below the Method Detection Li
N/A - Not Applicable

N/A - Not Applicable (a) - Data reported in ug/kg.

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Client Comple ID			01BS-0-2	01BS-10-12	01BS-10-12 01BS-14-16 02B-S-0-2 02B-S-12-14 02B-S-18-20	02B-S-0-2	02B-S-12-14	02B-S-18-20	03BS-0-2	03BS-6-8
I ob Somple Number			29017005	29017006	29017007	28934001	28934002	28934003	29017011	29017012
Motrie			SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL
Somple Date			3/6/91	3/6/91	3/6/91	2/28/91	2/28/91	2/28/91	3/6/91	3/6/91
Sample Date	CRDL	Action							South College	
Semivolatile Compounds	ug/kg	Levels (ug/kg)								
N-Nitroso-di-n-propylamine	330	100								
Hexachloroethane	330	80,000								
Isophorone	330	2E+06								
2-Nitrophenol	330	N/A								
2.4-Dimethylphenol	330	N/A								
his(2-Chloroethoxy)Methane	330	N/A								
2.4-Dichlorophenol	330	200,000								
1.2.4-Trichlorobenzene	330	2E+06								
4-Chloroaniline	330	N/A								
Hexachlorobutadiene	330	000'06								

Notes:

B - Applies to organic data only. Present in the corresponding method blank

J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.

E - Applies to Gas Chromatography/Mass Spectroscopy data only. Indicate compound above or below linear range of instrument. CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a)

Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

DANNELLY ANG - Montgomery, Alabama (a)

Client Sample ID			01BS-0-2	01BS-10-12	01BS-14-16	02B-S-0-2	01BS-10-12 01BS-14-16 02B-S-0-2 02B-S-12-14 02B-S-18-20	02B-S-18-20	03BS-0-2	03BS-6-8
Lab Sample Number			29017005	29017006	29017007	28934001	28934002	28934003	29017011	29017012
Matrix			SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL
Sample Date			3/6/91	3/6/91	3/6/91	2/28/91	2/28/91	2/28/91	3/6/91	3/6/91
	CRDL	Action	が重要を含むない		三 新音 表 卷 安全					医三角性
Semivolatile Compounds	ug/kg	Levels (ug/kg)		经过的效应					金属的有限的特殊之	· 有一种 · · · · · · · · · · · · · · · · · · ·
4-Chloro-3-methylphenol	330	N/A			-					
Hexachlorocyclopentadiene	330	000,009								
2,4,6-Trichlorophenol	330	40,000								
2,4,5-Trichlorophenol	1600	8E+06								
2-Chloronaphthalene	330	N/A								
2-Nitroaniline	1600	N/A								
Dimethyl Phthalate	330	N/A								
Acenaphthylene	330	N/A								
2,6-Dinitrotoluene	330	1,000								
3-Nitroaniline	1600	N/A								

Notes:

- B Applies to organic data only. Present in the corresponding method blank
- J Applies to organic data only. Value detected is greater than zero but less than the CRDL. E Applies to Gas Chromatography/Mass Spectroscopy data only.
 - Applies to Gas Chromatography/Mass Spectroscopy data only.
 Indicate compound above or below linear range of instrument.

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254,521(a)

Actual Egyesis proposed in Appendix A of 40C1 N234:221(a)
Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

DANNELLY ANG - Montgomery, Alabama (*) SITE 2, OIL WATER SEPARATOR 187th Fighter Group

Client Sample ID			01BS-0-2	01BS-10-12	01BS-14-16	02B-S-0-2	02B-S-12-14	02B-S-18-20	03BS-0-2	03BS-6-8
I ob Somple Number			29017005	29017006	29017007	28934001	28934002	28934003	29017011	29017012
Motrix			SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	TIOS	TIOS
Sample Date			3/6/91	3/6/91	3/6/91	2/28/91	2/28/91	2/28/91	3/6/91	3/6/91
	CRDL	Action								
Semivolatile Compounds	ug/kg	Levels (ug/kg)								
Acenaphthene	330	N/A								
2.4-Dintrophenol	1600	200,000								
4-Nitrophenol	1600	N/A								
Dibenzofuran	330	N/A								
2,4-Dinitrotoluene	330	N/A								
Diethylphthalate	330	6E+07								
4-Chlorophenyl-phenylether	330	N/A								
4-Nitroaniline	1600	N/A								
4,6-Dinitro-2-methylphenol	1600	N/A								
4-Bromophenyl-phenylether	330	N/A								

B - Applies to organic data only. Present in the corresponding method blank J - Applies to organic data only. Value detected is greater than zero but less than the CRDL. E - Applies to Gas Chromatography/Mass Spectroscopy data only.

Indicate compound above or below linear range of instrument.

CRDL - Contract Required Detection Limit

Blank Space - Value is below the Method Detection Limit(MDL). Action Levels proposed in Appendix A of 40CFR254.521(a)

N/A - Not Applicable

DANNELLY ANG - Montgomery, Alabama (*)

Client Sample ID			01BS-0-2	01BS-10-12	01BS-14-16		02B-S-12-14	02B-S-18-20	03BS-0-2	03BS-6-
Lab Sample Number			29017005	29017006	29017007	2	28934002 28934003	28934003	29017011	2901701
Matrix			SOIL	SOIL	TIOS	TIOS	SOIL	SOIL	SOIL	SOIL
Sample Date			3/6/91	3/6/91	16/9/8	2/28/91	2/28/91	2/28/91	3/6/91	3/6/91
计多数形式 网络多种 医神经囊炎 的复数	CRDL	Action		计算数据						
Semivolatile Compounds	ug/kg	Levels (ug/kg)								
Hexachlorobenzene	330	N/A								
Pentachlorophenol	1600	2E+06								
Butylbenzylphthalate	330	2E+07								
3,3'-Dichlorobenzidine	099	2,000								
Di-n-octylphthalate	330	N/A	-							

Notes:

- B Applies to organic data only. Present in the corresponding method blank
- J Applies to organic data only. Value detected is greater than zero but less than the CRDL.
 - E Applies to Gas Chromatography/Mass Spectroscopy data only.

Indicate compound above or below linear range of instrument.

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a)

Blank Space - Value is below the Method Detection Limit(MDL). N/A - Not Applicable

DANNELLY ANG - Montgomery, Alabama (*)

Client Sample ID			03BS-12-14	03BS-12-14D	03BS12-14 MS	03BS12-14 MS 03BS12-14 DUP	04BS-0-2	04BS-6-8	04BS-14-16
Lab Sample Number			29017013	29017014	29017M14	29017P14	29017001	29017002	29017003
Matrix			SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL
Sample Date			3/6/91	3/6/91	3/6/91	3/6/91	3/6/91	3/6/91	3/6/91
(の) かんこう はいかい ないない ないかい ないない ないかい ないかい ないかい ないかい	CRDL	Action						· · · · · · · · · · · · · · · · · · ·	このなるとなると
Semivolatile Compounds	ug/kg	Levels (ug/kg)							
2-Methylnaphthalene	330	N/A	-				4100	43 J	170 J
Anthracene	330	N/A							
Benzo(a)anthracene	330	N/A							
Benzo(a)pyrene	330	N/A							
Benzo(b)fluoranthene	330	N/A							
Benzo(g,h,i)perylene	330	N/A							
Benzo(k)fluoranthene	330	N/A							-
Benzoic Acid	1600	N/A							
Chrysene	330	N/A							
Di-n-butylphthalate	330	8E+06					-	50 J	48 J

Notes:

B - Applies to organic data only. Present in the corresponding method blank

J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.

E - Applies to Gas Chromatography/Mass Spectroscopy data only. Indicate compound above or below linear range of instrument.

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a) Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

DANNELLY ANG - Montgomery, Alabama (*)

Client Sample ID			03BS-12-14	03BS-12-14D	03BS12-14 MS	03BS12-14 DUP	04BS-0-2	04BS-6-8	04BS-14-16
Lab Sample Number			29017013	29017014	29017M14	29017P14	29017001	.29017002	29017003
Matrix			SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOLL
Sample Date			3/6/91	3/6/91	3/6/91	3/6/91	3/6/91	3/6/91	3/6/91
	CRDL	Action							
Semivolatile Compounds	ug/kg	Levels (ug/kg)							
Dibenz(a,h)anthracene	330	N/A							
Fluoranthene	330	N/A							
Fluorene	330	N/A					100 J		·
Indeno(1,2,3-cd)pyrene	330	N/A					,		
N-Nitrosodiphenylamine	330	100,000					94 J		
Naphthalene	330	N/A	48 J	48 J			8800 E	60 J	340 J
Nitrobenzene	330	40,000					270 J		
Phenanthrene	330	N/A					110 J		
Pyrene	330	N/A					44 J		
Bis(2-Ethylhexyl)phthalate	330	50,000	270 BJ	320 BJ			540 B	820 B	1500 B

Notes:

B - Applies to organic data only. Present in the corresponding method blank J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.

E - Applies to Gas Chromatography/Mass Spectroscopy data only.
 Indicate compound above or below linear range of instrument.

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a)

Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

DANNELLY ANG - Montgomery, Alabama "

Client Semale ID			03BS-12-14	03BS-12-14D	03BS12-14 MS	03BS12-14 DUP	04BS-0-2	04BS-6-8	04BS-14-16
I ob Somple Number			29017013	29017014	29017M14	29017P14	29017001	29017002	29017003
Motrix			SOIL	SOIL	SOIL	SOIL	SOIL	SOL	SOIL
Sample Date			3/6/91	3/6/91	3/6/91	3/6/91	3/6/91	3/6/91	3/6/91
は、これでは、一般のでは、一般のでは、一般のでは、一般のでは、一般のでは、一般のでは、一般のでは、一般のでは、一般のでは、一般のでは、一般のでは、一般のでは、一般のでは、一般のでは、一般のでは、一般のでは、	CRDL	Action							· · · · · · · · · · · · · · · · · · ·
Semivolatile Compounds	ug/kg	Levels (ug/kg)							· · · · · · · · · · · · · · · · · · ·
Phenol	330	SE+07							
bis(2-Chloroethyl)Ether	330	N/A							
2-Chlorophenol	330	400,000							
1,3-Dichlorobenzene	330	N/A			-				
1.4-Dichlorobenzene	330	N/A							
Benzyl Alcohol	330	N/A							
1.2-Dichlorobenzene	330	N/A							
2-Methylphenol	330	N/A							
bis(2-Chloroisopropyl)Ether	330	N/A							
4-Methylphenol	330	N/A							

Notes

B - Applies to organic data only. Present in the corresponding method blank

J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.

E - Applies to Gas Chromatography/Mass Spectroscopy data only.

Indicate compound above or below linear range of instrument.

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a)

Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

DANNELLY ANG - Montgomery, Alabama (*)

Client Sample ID			03BS-12-14	03BS-12-14D	03BS12-14 MS	03BS12-14 MS 03BS12-14 DUP	04BS-0-2	04BS-6-8	04BS-14-16
Lab Sample Number			29017013	29017014	29017M14	29017P14	29017001	29017002	29017003
Matrix			TIOS	SOIL	SOIL	TIOS	SOIL	SOL	SOIL
Sample Date			3/6/91	3/6/91	3/6/91	16/9/8	3/6/91	3/6/91	3/6/91
	CRDL	Action	と言うのではない。	《 100 100 100 100 100 100 100 100 100 10	· · · · · · · · · · · · · · · · · · ·				
Semivolatile Compounds	ug/kg	Levels (ug/kg)	· 关键的 1000 1000 1000 1000 1000 1000 1000 10						· 工作 专门的警卫的职业第
N-Nitroso-di-n-propylamine	330	100							
Hexachloroethane	330	80,000							
Isophorone	330	2E+06							
2-Nitrophenol	330	N/A							
2,4-Dimethylphenol	330	N/A							
bis(2-Chloroethoxy)Methane	330	N/A							
2,4-Dichlorophenol	330	200,000							
1,2,4-Trichlorobenzene	330	2E+06	-						
4-Chloroaniline	330	N/A							
Hexachlorobutadiene	330	90,000							

Notes:

B - Applies to organic data only. Present in the corresponding method blank

J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.

E - Applies to Gas Chromatography/Mass Spectroscopy data only. Indicate compound above or below linear range of instrument.

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a)

Blank Space - Value is below the Method Detection Limit(MDL). N/A - Not Applicable

DANNELLY ANG - Montgomery, Alabama 🙉 SITE 2, OIL WATER SEPARATOR 187th Fighter Group

Client Sample ID			03BS-12-14	03BS-12-14D	03BS12-14 MS	03BS12-14 DUP	04BS-0-2	04BS-6-8	04BS-14-16
Lab Sample Number			29017013	29017014	29017M14	29017P14	29017001	29017002	29017003
Matrix			SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL
Sample Date			3/6/91	3/6/91	3/6/91	3/6/91	3/6/91	3/6/91	3/6/91
いい まからをなっている あれるから	CRDL	Action					のこととの意味の対象	The second of th	
Semivolatile Compounds	ug/kg	Levels (ug/kg)					Selection and the second	· · · · · · · · · · · · · · · · · · ·	
4-Chloro-3-methylphenol	330	N/A							
Hexachlorocyclopentadiene	330	000'009							
2,4,6-Trichlorophenol	330	40,000							
2,4,5-Trichlorophenol	1600	8E+06							
2-Chloronaphthalene	330	N/A							
2-Nitroaniline	1600	N/A							
Dimethyl Phthalate	330	N/A							
Acenaphthylene	330	N/A							
2,6-Dinitrotoluene	330	1,000							
3-Nitroaniline	1600	N/A							

Notes:

B - Applies to organic data only. Present in the corresponding method blank J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.

E - Applies to Gas Chromatography/Mass Spectroscopy data only.

Indicate compound above or below linear range of instrument.

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a)

Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable (a) - Data reported in ug/kg.

DANNELLY ANG - Montgomery, Alabama (*)

Client Sample ID			03BS-12-14	03BS-12-14D	03BS12-14 MS	03BS12-14 MS 03BS12-14 DUP	04BS-0-2	04BS-6-8	04BS-14-16
Lab Sample Number			29017013	29017014	29017M14	29017P14	29017001	29017002	29017003
Matrix			SOIL	TIOS	SOIL	SOIL	SOIL	SOIL	SOIL
Sample Date			3/6/91	16/9/8	3/6/91	3/6/91	3/6/91	3/6/91	3/6/91
	CRDL	Action		· 特别,以第二次的			。 一种性學學學學學學學學學學學學學學學學學學學學學學學學學學學學學學學學學學學學		
Semivolatile Compounds	ug/kg	Levels (ug/kg)		计型机器 非常教科教					新文章(1) \$P\$ (1) \$P\$ (1) \$P\$ (1)
Acenaphthene	330	N/A							
2,4-Dintrophenol	1600	200,000							
4-Nitrophenol	1600	N/A							
Dibenzofuran	330	N/A							
2,4-Dinitrotoluene	330	N/A							
Diethylphthalate	330	6E+07							
4-Chlorophenyl-phenylether	330	N/A							
4-Nitroaniline	1600	N/A							
4,6-Dinitro-2-methylphenol	1600	N/A							
4-Bromophenyl-phenylether	330	N/A							

Notes:

- B Applies to organic data only. Present in the corresponding method blank J Applies to organic data only. Value detected is greater than zero but less than the CRDL.
- E Applies to Gas Chromatography/Mass Spectroscopy data only.

Indicate compound above or below linear range of instrument.

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a)

Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

DANNELLY ANG - Montgomery, Alabama 🗝 SITE 2, OIL WATER SEPARATOR 187th Fighter Group

Client Semple ID			03BS-12-14	03BS-12-14D	03BS12-14 MS	03BS12-14 MS 03BS12-14 DUP	04BS-0-2	04BS-6-8	04BS-14-16
Lab Sample Number			29017013	29017014	29017M14	29017P14	29017001	29017002	29017003
Matrix			SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL
Sample Date			3/6/91	3/6/91	3/6/91	3/6/91	3/6/91	3/6/91	3/6/91
	CRDL	Action							
Semivolatile Compounds	ug/kg	Levels (ug/kg)			×				
Hexachlorobenzene	330	N/A							
Pentachlorophenol	1600	2E+06							
Butylbenzylphthalate	330	2E+07		·					
3,3'-Dichlorobenzidine	099	2,000							
Di-n-octylphthalate	330	N/A							

Notes:

B - Applies to organic data only. Present in the corresponding method blank

J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.

E - Applies to Gas Chromatography/Mass Spectroscopy data only.

Indicate compound above or below linear range of instrument. CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a) Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

DANNELLY ANG - Montgomery, Alabama (*)

Client Sample ID			04BS-14-16D	05BS-0-2	05BS-8-10	05BS-12-14
Lab Sample Number			29017004	29017008	29017009	29017010
Matrix			TIOS	TIOS	TIOS	SOIL
Sample Date			3/6/91	3/6/91	3/6/91	3/6/91
	CRDL	Action				
Semivolatile Compounds	ug/kg	Levels (ug/kg)	productive makes			
2-Methylnaphthalene	330	N/A	180 J		1200	69 J
Anthracene	330	N/A				
Benzo(a)anthracene	330	N/A			52 J	
Benzo(a)pyrene	330	N/A			<i>57</i> J	
Benzo(b)fluoranthene	330	N/A			83 J	
Benzo(g,h,i)perylene	330	N/A			50 J	
Benzo(k)fluoranthene	330	N/A			64 J	
Benzoic Acid	1600	N/A				780 J
Chrysene	330	N/A			84 J	
Di-n-butylphthalate	330	8E+06				

B - Applies to organic data only. Present in the corresponding method blank J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.

E - Applies to Gas Chromatography/Mass Spectroscopy data only.

Indicate compound above or below linear range of instrument.

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a)

Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

DANNELLY ANG - Montgomery, Alabama (*) SITE 2, OIL WATER SEPARATOR 187th Fighter Group

Client Sample ID			04BS-14-16D	05BS-0-2	05BS-8-10	05BS-12-14
Lab Sample Number			29017004	29017008	29017009	29017010
Matrix			SOIL	SOIL	SOIL	SOIL
Sample Date			3/6/91	3/6/91	3/6/91	3/6/91
変更を持ちますが、カンラングでは、おからからできた。	CRDL	Action				
Semivolatile Compounds	ug/kg	Levels (ug/kg)				
Dibenz(a,h)anthracene	330	N/A				
Fluoranthene	330	N/A			92 J	
Fluorene	330	N/A				
Indeno(1,2,3-cd)pyrene	330	N/A			<i>57</i> J	
N-Nitrosodiphenylamine	330	100,000	<i>S7</i> J		82 J	
Naphthalene	330	N/A	420		1100	92 J
Nitrobenzene	330	40,000				
Phenanthrene	330	N/A			49 J	
Pyrene	330	N/A			67 J	
Bis(2-Ethylhexyl)phthalate	330	50,000	2300 B	1000 B	6100 B	2700 B

B - Applies to organic data only. Present in the corresponding method blank J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.

E - Applies to Gas Chromatography/Mass Spectroscopy data only. Indicate compound above or below linear range of instrument.

CRDL - Contract Required Detection Limit

Blank Space - Value is below the Method Detection Limit(MDL). Action Levels proposed in Appendix A of 40CFR254.521(a)

N/A - Not Applicable (a) - Data reported in ug/kg.

DANNELLY ANG - Montgomery, Alabama (*)

Client Sample ID			04BS-14-16D	05BS-0-2	05BS-8-10	05BS-12-14
Lab Sample Number		-	29017004	29017008	29017009	29017010
Matrix			TIOS	SOIL	TIOS	SOIL
Sample Date			3/6/91	3/6/91	3/6/91	3/6/91
是一个一个一个一个一个一个一个一个一个一个一个一个一个一个一个一个一个一个一个	CRDL	Action				
Semivolatile Compounds	ug/kg	Levels (ug/kg)	非国际的国际的国际的国际			
Phenol	330	5E+07				
bis(2-Chloroethyl)Ether	330	N/A				
2-Chlorophenol	330	400,000				
1,3-Dichlorobenzene	330	N/A				
1,4-Dichlorobenzene	330	N/A				
Benzyl Alcohol	330	N/A				
1,2-Dichlorobenzene	330	N/A				
2-Methylphenol	330	N/A				
bis(2-Chloroisopropyl)Ether	330	N/A				
4-Methylphenol	330	N/A				

B - Applies to organic data only. Present in the corresponding method blank

J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.

E - Applies to Gas Chromatography/Mass Spectroscopy data only.

Indicate compound above or below linear range of instrument. CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a)

Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

DANNELLY ANG - Montgomery, Alabama (*)

Client Sample ID			04BS-14-16D	05BS-0-2	05BS-8-10	05BS-12-14
Lab Sample Number			29017004	29017008	29017009	29017010
Matrix			SOIL	SOIL	SOIL	SOIL
Sample Date			3/6/91	3/6/91	3/6/91	3/6/91
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Semivolatile Compounds	ug/kg	Levels (ug/kg)				
N-Nitroso-di-n-propylamine	330	100				
Hexachloroethane	330	80,000				
Isophorone	330	2E+06				
2-Nitrophenol	330	N/A				
2,4-Dimethylphenol	330	N/A				
bis(2-Chloroethoxy)Methane	330	N/A				
2,4-Dichlorophenol	330	200,000				
1,2,4-Trichlorobenzene	330	2E+06				
4-Chloroaniline	330	N/A				
Hexachlorobutadiene	330	90,000				

Notes:

B - Applies to organic data only. Present in the corresponding method blank

J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.

E - Applies to Gas Chromatography/Mass Spectroscopy data only.

Indicate compound above or below linear range of instrument.

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a)

Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

DANNELLY ANG - Montgomery, Alabama (a)

Client Sample ID			04BS-14-16D	05BS-0-2	05BS-8-10	05BS-12-14
Lab Sample Number			29017004	29017008	29017009	29017010
Matrix			TIOS	TIOS	SOIL	SOIL
Sample Date			3/6/91	3/6/91	3/6/91	3/6/91
	CRDL	Action				
Semivolatile Compounds	ug/kg	Levels (ug/kg)	们是有 从 。但在4月14年的			
4-Chloro-3-methylphenol	330	N/A				
Hexachlorocyclopentadiene	330	000,009				
2,4,6-Trichlorophenol	330	40,000				
2,4,5-Trichlorophenol	1600	8E+06	,			
2-Chloronaphthalene	330	N/A				
2-Nitroaniline	1600	N/A				
Dimethyl Phthalate	330	N/A				
Acenaphthylene	330	N/A				
2,6-Dinitrotoluene	330	1,000				
3-Nitroaniline	1600	N/A				

Notes:

- B Applies to organic data only. Present in the corresponding method blank
- J Applies to organic data only. Value detected is greater than zero but less than the CRDL.
 - E Applies to Gas Chromatography/Mass Spectroscopy data only.

Indicate compound above or below linear range of instrument.

DIV - Contract Demitsed Detection I imit

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a) Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

DANNELLY ANG - Montgomery, Alabama (*)

Client Sample ID			04BS-14-16D	05BS-0-2	05BS-8-10	05BS-12-14
Lab Sample Number			29017004	29017008	29017009	29017010
Matrix			SOIL	SOIL	TIOS	SOIL
Sample Date			3/6/91	3/6/91	3/6/91	3/6/91
できる。 第二章を表する。 多生がある。 ない	CRDL	Action	。 《四十二章》			
Semivolatile Compounds	ug/kg	Levels (ug/kg)				
Acenaphthene	330	N/A				
2,4-Dintrophenol	1600	200,000	,			
4-Nitrophenol	1600	N/A				
Dibenzofuran	330	N/A				
2,4-Dinitrotoluene	330	N/A				
Diethylphthalate	330	6E+07				
4-Chlorophenyl-phenylether	330	N/A			٠	
4-Nitroaniline	1600	N/A				
4,6-Dinitro-2-methylphenol	1600	N/A				
4-Bromophenyl-phenylether	330	N/A				

Notes:

B - Applies to organic data only. Present in the corresponding method blank

J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.

E - Applies to Gas Chromatography/Mass Spectroscopy data only.

Indicate compound above or below linear range of instrument.

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a)

Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

DANNELLY ANG - Montgomery, Alabama (*)

Client Sample ID			04BS-14-16D	05BS-0-2	05BS-8-10	05BS-12-14
Lab Sample Number			29017004	29017008	29017009	29017010
Matrix			SOIL	TIOS	TIOS	SOIL
Sample Date			3/6/91	3/6/91	16/9/8	3/6/91
	CRDL	Action				
Semivolatile Compounds	ug/kg	Levels (ug/kg)				
Hexachlorobenzene	330	N/A				
Pentachlorophenol	1600	2E+06				
Butylbenzylphthalate	330	2E+07				
3,3'-Dichlorobenzidine	099	2,000				
Di-n-octylphthalate	330	N/A				

Notes:

B - Applies to organic data only. Present in the corresponding method blank

J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.

E - Applies to Gas Chromatography/Mass Spectroscopy data only. Indicate compound above or below linear range of instrument.

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a)

Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

DANNELLY ANG - Montgomery, Alabama (*)

Client Sample ID			01BS-0-2	01BS-10-12	01BS-14-16	02B-S-0-2	02B-S-12-14	02B-S-18-20	03BS-0-2	03BS-6-8
Lab Sample Number			29017005	29017006	29017007	28934001	28934002	28934003	29017011	29017012
Matrix			SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL
Sample Date			3/6/91	3/6/91	3/6/91	2/28/91	2/28/91	2/28/91	3/6/91	3/6/91
	CRDL	Action								
Inorganics	mg/kg	Levels (mg/kg)								
Antimony	16.2	30,000			5.3 B			3.7 B		
Arsenic	2.7	80,000			5.1			7.6		
Barium	54	4E+06			33.6 B			87.1		
Cadmium	1.4	40,000			0.81 B			0.75 B		
Chromium	2.7	400,000			19.0			21.5		
Copper	6.8	400,000			37.4			13.8		
Lead	0.8	N/A			5.6			7.8		
Mercury	0.05	20,000								
Nickel	10.8	200,000			25.1			19.8		
Selenium	1.4	N/A			0.71B					
Silver	2.7	200,000								
Thallium	2.7	000'9			0.28 B					
Zinc	5.4	4E+06			65.6			53.5		

Notes:

B - Value detected is less than the CRDL but greater than of equal to the MDL.

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a)

Blank Space - Value is below the Method Detection Limit(MDL).

CRDL expressed in mg/kg = {CRDL in ug/L * Digestion Factor (Assuming 200) * Unit Conversion(1/1000)}/%Solids(Assuming 75%)

N/A - Not Applicable

N/C - No change in the RPD.

(c) - Pelative Percent Difference(RPD) for duplicate samples.

(d) - Percent Recovery (%R) for Matrix Spike(MS) samples.(a) - Data reported in mg/kg.

DANNELLY ANG - Montgomery, Alabama (*)

14 03BS-12-14D 03BS12-14 MS 03BS12-14 DUP 0 3 29017014 29017M14 29017P14 2	03BS-12-14D 03BS12-14 MS 03BS12-14 DUP 29017014 29017M14 29017P14	03BS12-14 MS 03BS12-14 DUP 29017M14 29017P14		04BS-0- 2901700	12	04BS-6-8 29017002	04BS-14-16 29017003	04BS-14-16D 29017004
SOIL SOIL (d) SOIL(c)	SOIL SOIL (d)	SOIL (d)	SOIL (c)		SOIL	SOIL	SOIL	
3/6/91 3/6/91 3/6/91	3/6/91 3/6/91	3/6/91	3/6/91		3/6/91	3/6/91	3/6/91	
CRDL Action Action				***				黛娜
			金剛 李素明 排出					
16.2 30,000 4.1B 3.1 B 62.4 59.7	3.1 B 62.4	62.4	59.7			5.2 B		
80,000 91.8 N/C			N/C					
4E+06 31.3 B 50.3 B 92.1 8.0	50.3 B 92.1	92.1	8.0	1		55.7		
40,000 0.65 B 0.48 B 81.0 38.7	0.48 B 81.0	81.0	38.7			0.45 B		
171.4	17.8 171.4	171.4	35.4			16.8		
400,000 18.9 18.2 91.2 0.9	18.2 91.2	91.2	0.9			11.3		-
	6.3 106.7	106.7	3.3	\vdash		5.2		
0.05 20,000 N/C			N/C					
10.8 200,000 16.3 14.6 88.9 12.7	14.6 88.9	88.9	12.7			15.4		
N/A 0.97 B 3.0 41.9 23.7	3.0 41.9	41.9	23.7			0.40 B		
200,000 85.2 N/C			N/C					
6,000 0.23 B 84.4 N/C	84.4	84.4	N/C					
4E+06 39.6 49.0 92.6 11.3	49.0 92.6	92.6	11.3			48.0		

Notes:

B - Value detected is less than the CRDL but greater than of equal to the MDL.

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a)

Blank Space - Value is below the Method Detection Limit(MDL).

CRDL expressed in mg/kg = (CRDL in ug/L * Digestion Factor (Assuming 200)

* Unit Conversion(1/1000)}/%Solids(Assuming 75%)

N/A - Not Applicable

N/C - No change in the RPD.

(c) - Pelative Percent Difference(RPD) for duplicate samples. (d) - Percent Recovery (%R) for Matrix Spike(MS) samples. (a) - Data reported in mg/kg.

DANNELLY ANG - Montgomery, Alabama (*)

B - Value detected is less than the CRDL but greater than of equal to the MDL.

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a)

Blank Space - Value is below the Method Detection Limit(MDL).

CRDL expressed in mg/kg = {CRDL in ug/L * Digestion Factor (Assuming 200)

* Unit Conversion(1/1000)}/%Solids(Assuming 75%)

N/A - Not Applicable

N/C - No change in the RPD.

(c) - Pelative Percent Difference(RPD) for duplicate samples.

(d) - Percent Recovery (%R) for Matrix Spike(MS) samples.(a) - Data reported in mg/kg.

SITE 2, OIL WATER SEPARATOR 187th Fighter Group DANNELLY ANG - Montgomery, Alabama (a)

_	_		_			_	_	- 55	_	102.06	_	44.4	439
O5BS	17978	SOIL	3/6/91		A STATE OF THE STA		5.5		5.7		12.1		
O4BSDUP**	17978	SOIL	3/6/91	通过的编码的编码								9.9	
O4BS**	17978	SOIL	3/6/91				2120	69.1				< 1.1	
O4BSDUP	17947	SOIL	3/1/91									113	
O4BS	17947	SOIL	3/1/91				12.8	95.1	The second secon			378	
O3BSDUP	17978	SOIL	3/6/91				6.0	The second secon		and the second s	< 1.1	the second of th	
O3BS	17978	SOIL	3/6/91	方の表示である。 第15章 (15章 15章 15章 15章 15章 15章 15章 15章 15章 15章			< 1.1	<1.1	Carlos Santas Carlos Santas Sa	The second secon	< 1.1	e de la Caración de Caración d	
OZBS	17946	SOIL	2/28/91				16	And the second s			<1.1		< 1.1
OIBS	17978	SOIL	3/6/91				<1.1			< 1.1		< 1.1	
Client Sample ID	Lab Sample Number	Matrix	Sample Date		Total Petroleum Hydrocarbons	Sample Depth(Feet)	0-2	8-9	8-10	10-12	12-14	14-16	18-20

Notes:

** - Indicated the boring was redrilled and sampled. Shaded areas indicate depth intervals not sampled.

SITE 4 (APRON SPILL AREA) 187th Fighter Group DANNELLY ANG - Montgomery, Alabama (*)

2.			410000	A 1 B S 4 10	A1BC-16-18	A1-88-0-6	A1-SS-6-24	A2BS-4-6	A2BS-12-14	A2BS-12-14 A2BS-12-14DUP
Citent Sample 1D			29032004	29032005	29032006	17894001	17894002	29046001	29046002	29046003
Motrix			SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL
Somele Dote			3/7/91	3/7/91	3/7/91	2/21/91	2/21/91	3/8/91	3/8/91	3/8/91
Sample Date	CPNI	Action					· · · · · · · · · · · · · · · · · · ·		A TOP MAN TO SHEET	
PAH Compounds	ng/kg	Levels (ug/kg)							意の意味が発展する	《沙埃多林河縣][標本編書
Anthracene	330	N/A								
Benzo(a)anthracene	330	N/A								
Benzo(a)pyrene	330	N/A								
Benzo(b)fluoranthene	330	N/A								
Benzo(g.h.i)perylene	330	N/A			·				,	
Benzo(k)fluoranthene	330	N/A								
Chrysene	330	N/A								
Dibenzo(a,h)anthracene	330	N/A								
Fluoranthene	330	N/A				37 JX				
Fluorene	330	N/A								
Indeno(1,2,3-cd)pyrene	330	N/A								
Phenanthrene	330	N/A								
Pyrene	330	N/A				38 JX				
Naphthalene	330	N/A								
2-Methylnaphthalene	330	N/A								
1-Methylnaphthalene	330	N/A								
Acenaphthylene	330	N/A								
Acenaphthene	330	N/A								

Notes

JX - The compound was detected and quantitated below CRDL. CRDL - Contract Required Detection Limit Action Levels proposed in Appendix A of 40CFR254.521(a) Blank Space - Value is below the Method Detection Limit(MDL). N/A - Not Applicable (a) - Data reported in ug/kg.

69-O

DANNELLY ANG - Montgomery, Alabama " SITE 4 (APRON SPILL AREA) 187th Fighter Group

Client Sample ID			A2BS-18-20	A2-SS-0-6	A2-SS-6-24	A3BS-0-2	A3BS-8-10	A3BS-18-20	A3BS-18-20D	A3-SS-0-6
Lab Sample Number			29046004	17894003	17894004	29032007	29032008	29032009	29032010	17894005
Matrix			SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL
Sample Date			3/8/91	2/21/91	2/21/91	3/7/91	3/7/91	3/7/91	3/7/91	2/21/91
	CRDL	Action			· · · · · · · · · · · · · · · · · · ·					
PAH Compounds	ug/kg	Levels (ug/kg)								
Anthracene	330	N/A								
Benzo(a)anthracene	330	N/A		088	190					0009
Benzo(a)pyrene	330	N/A		3000	220					12000
Benzo(b)fluoranthene	330	N/A		1500	240					9500
Benzo(g,h,i)perylene	330	N/A		1700	250					8800
Benzo(k)fluoranthene	330	N/A		880	130					5000
Chrysene	330	N/A		1900	260					7900
Dibenzo(a,h)anthracene	330	N/A		390	39 JX					950
Fluoranthene	330	N/A		1900	340					7400
Fluorene	330	N/A					·			
Indeno(1,2,3-cd)pyrene	330	N/A		2200	260					10000
Phenanthrene	330	N/A		420	140					500 JX
Pyrene	330	N/A		1500	280					8600
Naphthalene	330	N/A								
2-Methylnaphthalene	330	N/A								
1-Methylnaphthalene	330	N/A								
Acenaphthylene	330	N/A								
Acenaphthene	330	N/A			-					

JX - The compound was detected and quantitated below CRDL. CRDL - Contract Required Detection Limit Action Levels proposed in Appendix A of 40CFR254.521(a) Blank Space - Value is below the Method Detection Limit(MDL). N/A - Not Applicable (a) - Data reported in ug/kg.

SITE 4 (APRON SPILL AREA) 187th Fighter Group DANNELLY ANG - Montgomery, Alabama (*)

Client Comple ID			A3-SS-0-6-DUP	A3-SS-6-24	A3-SS-6-24DUP	A4-SS-0-6	A4-SS-6-24	A5-SS-0-6	A5-SS-6-24	A6-SS-0-6
I oh Sample Number			17894006	17894007	17894008	17894009	17894010	17894011	17894012	17894013
Motriv			SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL
Somula Data			2/21/91	2/21/91	2/21/91	2/21/91	2/21/91	2/21/91	2/21/91	2/21/91
Sample Date	CBDI	Action								
PAH Communds	ug/kg	Levels (ug/kg)			是 1. 1. 1. 1. 1. 1. 1. 1. 1. 1. 1. 1. 1.	2. 工艺教育2. 八点				お生活を発行する
Anthracene	330	N/A	Propried to the propried to th					1000	920	
Benzo(a)anthracene	330	N/A	130	110		720	750	4100	3800	710
Benzo(a)nvrene	330	N/A	310	230		1200	1000	5000	4600	1400
Benzo(b)fluoranthene	330	N/A	330	260		1400	088	4100	3700	1500
Benzo(o h i)nerylene	330	N/A	029	420	45 JX	1500	940	3400	3000	1200
Banzo(b)fluoranthene	330	N/A	170	120		700	490	2200	2000	810
Chrysene	330	N/A	250	230		1400	096	5300	4800	1600
Dibenzo(a h)anthracene	330	NA	80	56		200 JX	140 JX	530 JX	780	220 JX
Fluoranthene	330	N/A	260	260	43 JX	2200	1600	9400	8400	2700
Fluorene	330	N/A							300 JX	
Indepo(1.2.3-cd)nyrene	330	N/A	570	390	89	1400	006	4100	3800	1300
Phenanthrene	330	N/A	79	91		830	840	3300	2900	910
Pyrene	330	N/A	230	240		1800	1400	7600	7100	2200
Naphthalene	330	N/A		·						
2-Methylnaphthalene	330	N/A		,						
1-Methylnaphthalene	330	N/A		٠						
Acenaphthylene	330	N/A								
Acenaphthene	330	N/A								

Notes:

JX - The compound was detected and quantitated below CRDL. CRDL - Contract Required Detection Limit Action Levels proposed in Appendix A of 40CFR254.521(a) Blank Space - Value is below the Method Detection Limit(MDL). N/A - Not Applicable (a) - Data reported in ug/kg.

SITE 4 (APRON SPILL AREA) 187th Fighter Group

DANNELLY ANG - Montgomery, Alabama (*)

			A6-SS-6-24	A7-SS-0-6	A7-SS-6-24	A8-SS-0-6	A8-SS-6-24
Lab Sample Number			17894014	17894015	17894016	17894017	17894018
Matrix			SOIL	SOIL	SOIL	SOIL	SOIL
Sample Date			2/21/91	2/22/91	2/22/91	2/22/91	2/22/91
	CRDL	Action		· 人民等的基础			
PAH Compounds	ug/kg	Levels (ug/kg)					
Anthracene	330	N/A					
Benzo(a)anthracene	330	N/A	4200	330 JX	068		
Benzo(a)pyrene	330	N/A	4500	640	1100		
Benzo(b)fluoranthene	330	N/A	4700	470	820		
Benzo(g,h,i)perylene	330	N/A	3600	1000	098		
Benzo(k)fluoranthene	330	N/A	2500	250 JX	480		
Chrysene	330	N/A	2900	069	1100		
Dibenzo(a,h)anthracene	330	N/A	700				
Fluoranthene	330	N/A	10000	280	1200	43 JX	
Fluorene	330	N/A					
Indeno(1,2,3-cd)pyrene	330	N/A	4200	1600	910		
Phenanthrene	330	N/A	4000		230 JX		
Pyrene	330	N/A	8300	260	1200	59 JX	
Naphthalene	330	N/A					
2-Methylnaphthalene	330	N/A		*			
1-Methylnaphthalene	330	N/A					
Acenaphthylene	330	N/A					
Acenaphthene	330	N/A					

Notes.

JX - The compound was detected and quantitated below CRDL. CRDL - Contract Required Detection Limit Action Levels proposed in Appendix A of 40CFR254.521(a) Blank Space - Value is below the Method Detection Limit(MDL). N/A - Not Applicable

DANNELLY ANG - Montgomery, Alabama (*) QC DATA 187th Fighter Group

Client Sample ID				TRAVEL BLANK TRAVEL BLANK TRAVEL BLANK TRAVEL BLANK TRAVEL BLANK TRAVEL BLANK TRIP BLANK TRV-02:2-28	TRIP BLANK	FRV-02-2-28					
Lab Sample Number				29032003	28973002	28998018	29046006	29017016	28909009	28891019	28934015
Matrix				WATER	WATER	WATER	WATER	WATER	WATER	WATER	WATER
Sample Date									2/26/91		2/28/91
からない 聖職は万里の世界には、大学のないか、	MCL	CRDL	Action								
Volatile Compounds	ng/L	ng/L	Levels (ug/L)								
trans-1,3-Dichloropropene	N/A	5	10								
4-Methyl-2-Pentanone	N/A	10	N/A								
2-Hexanone	N/A	10	N/A								
Tetrachloroethene	5	5	N/A								
1,1,2,2-Tetrachloroethane	N/A	5	2								
Toluene	1,000		10,000								
Chlorobenzene	N/A	5	200								
Ethyl Benzene	700	5	4,000								
Styrene	100	5	7,000								
Xylene (total)	10,000	5	7,000								

B - Applies to organic data only. Present in the corresponding method blank.

J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.

B - Applies to Gas Chromatography/Mass Spectroscopy data only.

Indicate compound above or below linear range of instrument.

Blank Space - Value is below the Method Detection Limit(MDL).

MCL - Maximum Contaminant Level (U.S. Drinking Water Standards)

CRDL - Contract Required Detection Limit
Action Levels proposed in Appendix A of 40CFR254.521(a)
N/A - Not Applicable
(a) - Data reported in ug/L.

QC DATA 187th Fighter Group

DANNELLY ANG - Montgomery, Alabama (*)

Client Sample ID				ERB-06-3-7	ERB-07-3-8	ER-1	FB-01-3-4	FB-02-3-7	FB-03-3-7
Lab Sample Number				29041001	29046005	28891020	28973001	29032002	29032001
Matrix	-			WATER	WATER	WATER	WATER	WATER	WATER
Sample Date				3/7/91	3/8/91		3/4/91	3/7/91	3/7/91
のできるというというできる。 のでは、これでは、これでは、これでは、これでは、これでは、これでは、これでは、これ	MCL	CRDL	Action						
Volatile Compounds	ug/L	ng/L	Levels (ug/L)		使物理性处理的增生				
Acetone	N/A	10	4,000					22 B	
Bromodichloromethane	100	5	0.03						14
Bromoform	100	5	700						4 J
Carbon Disulfide	N/A	5	4,000						
Chloroform	100	5	9		,				13
Dibromochloromethane	N/A	5	N/A						14
Methylene Chloride	N/A	5	5					6 B	4 BJ
Trichloroethene	5	5	N/A						
Chloromethane	N/A	10	N/A						
Bromomethane	N/A	10	50						
Vinyl Chloride	2	10	N/A						
Chloroethane	N/A	10	N/A						

Notes:

- B Applies to organic data only. Present in the corresponding method blank.
- J Applies to organic data only. Value detected is greater than zero but less than the CRDL.
 - E Applies to Gas Chromatography/Mass Spectroscopy data only.

Indicate compound above or below linear range of instrument.

Blank Space - Value is below the Method Detection Limit(MDL).

MCL - Maximum Contaminant Level (U.S. Drinking Water Standards)

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a)

N/A - Not Applicable

DANNELLY ANG - Montgomery, Alabama (a) 187th Fighter Group QC DATA

Client Sample ID				EQUIPMENT	EQUIPMENT ERB05-3-6 DUP ERB05-3-6 MS ERB-01-2-26 ERB-02-2-27 ERB-03-2-28 ERB-04-3-5 ERB-05-3-6	ERB05-3-6 MS	ERB-01-2-26	ERB-02-2-27	ERB-03-2-28	ERB-04-3-5	ERB-05-3-6
Lab Sample Number				17894020	29017P15	29017M15	28909010 28922004		28934014 28998017 29017015	28998017	29017015
Matrix				WATER	WATER	WATER	WATER	WATER	WATER	WATER	WATER
Sample Date					3/6/91	3/6/91	2/26/91	2/27/91	2/28/91	3/5/91	3/6/91
一 一 一 一 一 一 一 一 一 一 一 一 一 一 一 一 一 一 一	MCL	CRDL	Action								
Volatile Compounds	ng/L	ug/L	Levels (ug/L)								
Acetone	N/A	10	4,000				20 B	25 B			36 B
Bromodichloromethane	100	5	0.03								
Bromoform	100	5	700								
Carbon Disulfide	N/A	5	4,000								
Chloroform	100	5	9								
Dibromochloromethane	N/A	5	N/A								
Methylene Chloride	N/A	5	5				18 B	4 BJ			4 BJ
Trichloroethene	5	5	N/A								1 J
Chloromethane	N/A	10	N/A								
Bromomethane	N/A	10	50								
Vinyl Chloride	2	10	N/A								
Chloroethane	N/A	10	N/A								

B - Applies to organic data only. Present in the corresponding method blank.
J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.
E - Applies to Gas Chromatography/Mass Spectroscopy data only.
Indicate compound above or below linear range of instrument.

Blank Space - Value is below the Method Detection Limit(MDL).

MCL - Maximum Contaminant Level (U.S. Drinking Water Standards)

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a)

N/A - Not Applicable

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Client Sample ID				ERB-06-3-7	ERB-07-3-8	ER-1	FB-01-3-4	FB-02-3-7	FB-03-3-7	TRAVEL BLANK	ERB-06-3-7 ERB-07-3-8 ER-1 FB-01-3-4 FB-02-3-7 FB-03-3-7 TRAVEL BLANK TRAVEL BLANK
Lab Sample Number				29041001	29041001 29046005 28891020 28973001 29032002 29032001	28891020	28973001	29032002	29032001	29032003	28973002
Matrix				WATER	WATER WATER WATER WATER	WATER	WATER	WATER	WATER	WATER	WATER
Sample Date				377/91	3/8/91		3/4/91	377/91	377/91		
AND THE PROPERTY OF THE PROPER	MCL	CRDL	Action				1 9 THE	200			
Semivolatile Compounds	ng/L	ug/L	Levels (ug/L)								
Di-n-butylphthalate	N/A	10	N/A						33		
N-Nitrosodiphenylamine(1)	N/A	10	N/A								
Naphthalene	N/A	10	N/A					4 BJ			
bis(2-Ethylhexyl)phthalate	N/A	10	3					2 BJ	33	37	
Phenol	N/A	10	20,000								
bis(2-Chloroethyl)Ether	N/A	10	0.03								
2-Chlorophenol	N/A	10	200								
1,3-Dichlorobenzene	009	10	N/A								
1,4-Dichlorobenzene	750	10	N/A								
Benzyl Alcohol	N/A	10	N/A								

Notes:

B - Applies to organic data only. Present in the corresponding method blank.

J - Applies to organic data only. Value detected is greater than zero but Iess than the CRDL.

E - Applies to Gas Chromatography/Mass Spectroscopy data only.

Indicate compound above or below linear range of instrument.

MCL - Maximum Contaminant Level (U.S. Drinking Water Standards)

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a)

Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

(a) - Data reported in ug/L.

DANNELLY ANG - Montgomery, Alabama " QC DATA 187th Fighter Group

Client Samule ID				ERB-06-3-7	ERB-07-3-8	ER-1	FB-01-3-4	FB-02-3-7	FB-03-3-7	TRAVEL BLANK	ERB-06-3-7[ERB-07-3-8] ER-1 [FB-01-3-4 FB-02-3-7 FB-03-3-7 TRAVEL BLANK TRAVEL BLANK
I ab Sample Number				29041001	29041001 29046005 28891020 28973001 29032002 29032001	28891020	28973001	29032002	29032001	29032003	28973002
Matrix				WATER	WATER	WATER	WATER WATER WATER	WATER	WATER	WATER	WATER
Sample Date				377/91	3/8/91		3/4/91	3/7/91	3/7/91		
医外面的 医多种性 医多种性 医多种性 医多种性	MCL	CRDL	Action								生物位置國際宣言是接受的技术
S E	ug/L	ug/L	Levels (ug/L)							The state of t	以北京大学工艺学生 1000年 1000年 100
1,2-Dichlorobenzene	009	10	N/A								
2-Methylphenol	N/A	10	N/A								
bis(2-Chloroisopropyl)Ether	N/A	10	N/A								
4-Methylphenol	N/A	10	N/A								
N-Nitroso-di-n-propylamine	N/A	10	0.005								
Hexachloroethane	N/A	10	20								
Nitrobenzene	N/A	10	20								
Isophorone	N/A	10	06					.]			
2-Nitrophenol	N/A	10	N/A								
2,4-Dimethylphenol	N/A	50	N/A								

Notes:

B - Applies to organic data only. Present in the corresponding method blank.
J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.
E - Applies to Gas Chromatography/Mass Spectroscopy data only.
Indicate compound above or below linear range of instrument.
MCL - Maximum Contaminant Level (U.S. Drinking Water Standards)
CRDL - Contract Required Detection Limit
Action Levels proposed in Appendix A of 40CFR254,521(a)
Blank Space - Value is below the Method Detection Limit(MDL).
N/A - Not Applicable

DANNELLY ANG - Montgomery, Alabama (*)

Client Sample ID				ERB-06-3-7	ERB-07-3-8	ER-1	FB-01-3-4	FB-02-3-7	FB-03-3-7	TRAVEL BLANK	ERB-06-3-7 ERB-07-3-8 ER-1 FB-01-3-4 FB-02-3-7 FB-03-3-7 TRAVEL BLANK TRAVEL BLANK
Lab Sample Number				29041001	29041001 29046005 28891020 28973001 29032002 29032001	28891020	28973001	29032002	29032001	29032003	28973002
Matrix				WATER	WATER WATER WATER WATER	WATER	WATER	WATER	WATER	WATER	WATER
Sample Date				3/7/91	3/8/91		3/4/91	3/7/91	3/7/91		
	MCL	CRDL	Action								
Semivolatile Compounds	ng/L	ng/L	Levels (ug/L)		1001年間開始						が、一般の表現を表現
Benzoic Acid	N/A	10	N/A								
bis(2-Chloroethoxy)Methane	N/A	10	N/A								
2,4-Dichlorophenol	N/A	10	100								
1,2,4-Trichlorobenzene	0/	10	700								
4-Chloroaniline	N/A	10	N/A								
Hexachlorobutadiene	N/A	10	4								
4-Chloro-3-methylphenol	N/A	10	N/A								
2-Methylnaphthalene	N/A	10	N/A								
Hexachlorocyclopentadiene	N/A	10	200								
2,4,6-Trichlorophenol	N/A	10	3								

B - Applies to organic data only. Present in the corresponding method blank.
J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.
E - Applies to Gas Chromatography/Mass Spectroscopy data only.
Indicate compound above or below linear range of instrument.
MCL - Maximum Contaminant Level (U.S. Drinking Water Standards)

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a) Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

DANNELLY ANG - Montgomery, Alabama 🗝

Client Sample ID				ERB-06-3-7	ERB-07-3-8	ER-1	FB-01-3-4	FB-02-3-7	FB-03-3-7	TRAVEL BLANK	ERB-06-3-7 ERB-07-3-8 ER-1 FB-01-3-4 FB-02-3-7 FB-03-3-7 TRAVEL BLANK TRAVEL BLANK
Lab Sample Number				29041001	29041001 29046005 28891020 28973001 29032002 29032001	28891020	28973001	29032002	29032001	29032003	28973002
Matrix				WATER	WATER WATER WATER WATER	WATER	WATER	WATER	WATER	WATER	WATER
Sample Date				3/7/91	3/8/91		3/4/91	3/7/91	3/7/91		
· 1975年 · 1985年 · 198	MCL	CRDL	Action								
Semivolatile Compounds	ng/L	ug/L	Levels (ug/L)								
2,4,5-Trichlorophenol	N/A	50	4000								
2-Chloronaphthalene	N/A	10	N/A								
2-Nitroaniline	N/A	50	N/A								
Dimethyl Phthalate	N/A	10	N/A								
Acenaphthylene	N/A	10	N/A								
2,6-Dinitrotoluene	N/A	10	0.05								
3-Nitroaniline	N/A	50	N/A								
Acenaphthene	N/A	10	N/A								
2,4-Dintrophenol	N/A	50	70								
4-Nitrophenol	N/A	20	N/A								

B - Applies to organic data only. Present in the corresponding method blank.
 J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.
 E - Applies to Gas Chromatography/Mass Spectroscopy data only.

Indicate compound above or below linear range of instrument. MCL - Maximum Contaminant Level (U.S. Drinking Water Standards)

CRDL - Contract Required Detection Limit
Action Levels proposed in Appendix A of 40CFR254.521(a)
Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

DANNELLY ANG - Montgomery, Alabama (*)

Client Sample ID				ERB-06-3-7	ERB-07-3-8	ER-1	FB-01-3-4	FB-02-3-7	FB-03-3-7	TRAVEL BLANK	ERB-06-3-7 ERB-07-3-8 ER-1 FB-01-3-4 FB-02-3-7 FB-03-3-7 TRAVEL BLANK TRAVEL BLANK
Lab Sample Number				29041001	29041001 29046005 28891020 28973001 29032002 29032001	28891020	28973001	29032002	29032001	29032003	28973002
Matrix				WATER		WATER	WATER	WATER WATER WATER WATER	WATER	WATER	WATER
Sample Date				3/7/91	3/8/91		3/4/91	3/1/91	3/7/91		
是一位的 · · · · · · · · · · · · · · · · · · ·	MCL	CRDL	Action			THE STREET					
Semivolatile Compounds	ug/L	ng/L	Levels (ug/L)								が見り上が三個語彙
Dibenzofuran	N/A	10	N/A								
2,4-Dinitrotoluene	N/A	10	N/A								
Diethylphthalate	N/A	10	30,000								
4-Chlorophenyl-phenylether	N/A	10	N/A								
Fluorene	N/A	20	N/A								
4-Nitroaniline	N/A	20	N/A								
4,6-Dinitro-2-methylphenol	N/A	20	N/A								
4-Brómophenyl-phenylether	N/A	10	N/A								
Hexachlorobenzene	-	10	N/A								
Pentachlorophenol	1	20	1000								

B - Applies to organic data only. Present in the corresponding method blank. J - Applies to organic data only. Value detected is greater than zero but less than the CRDL. E - Applies to Gas Chromatography/Mass Spectroscopy data only.

MCL - Maximum Contaminant Level (U.S. Drinking Water Standards) Indicate compound above or below linear range of instrument.

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a) Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable (a) - Data reported in ug/L.

DANNELLY ANG - Montgomery, Alabama " QC DATA 187th Fighter Group

Client Comple ID				ERB-06-3-7	ERB-07-3-8	ER-1	FB-01-3-4	FB-02-3-7	FB-03-3-7	ERB-06-3-7 ERB-07-3-8 ER-1 FB-01-3-4 FB-02-3-7 FB-03-3-7 TRAVEL BLANK TRAVEL BLANK	TRAVEL BLANK
I ah Sample Number				29041001	29041001 29046005 28891020 28973001 29032002 29032001	28891020	28973001	29032002	29032001	29032003	28973002
Matrix				WATER	WATER	WATER WATER WATER WATER	WATER	WATER	WATER	WATER	WATER
Sample Date				3/7/91	3/8/91		3/4/91	3/7/91	3/7/91		
The second secon	MCL	CRDL	Action	新建設的	(A) (数) (1) (1)						
Semivolatile Compounds	ug/L	ng/L	Levels (ug/L)					2000年			
Phenanthrene	N/A	10	N/A								
Anthracene	N/A	10	N/A								
Fluoranthene	N/A	10	N/A								
Pyrene	N/A	10	N/A								
Butylbenzylphthalate	N/A	10	7,000								
3,3'-Dichlorobenzidine	N/A	20	0.08								
Benzo(a)anthracene	0.1	01	N/A								
Chrysene	0.2	10	N/A								
Di-n-octylphthalate	N/A	10	N/A								
Benzo(b)fluoranthene	0.2	10	N/A								

B - Applies to organic data only. Present in the corresponding method blank.

J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.

B - Applies to Gas Chromatography/Mass Spectroscopy data only.

Indicate compound above or below linear range of instrument.

MCL - Maximum Contaminant Level (U.S. Drinking Water Standards)

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a) Blank Space - Value is below the Method Detection Limit(MDL). N/A - Not Applicable (a) - Data reported in ug/L.

DANNELLY ANG - Montgomery, Alabama (*) 187th Fighter Group QC DATA

Client Sample ID				ERB-06-3-7	ERB-07-3-8	ER-1	FB-01-3-4	FB-02-3-7	FB-03-3-7	ERB-06-3-7 ERB-07-3-8 ER-1 FB-01-3-4 FB-02-3-7 FB-03-3-7 TRAVEL BLANK TRAVEL BLANK	TRAVEL BLANK
Lab Sample Number				29041001	29041001 29046005 28891020 28973001 29032002 29032001	28891020	28973001	29032002	29032001	29032003	28973002
Matrix				WATER	WATER WATER WATER WATER	WATER	WATER	WATER	WATER	WATER	WATER
Sample Date				3/7/91	16/8/8		3/4/91	3/7/91	377/91		
	MCL	CRDL	Action								
Semivolatile Compounds	ng/L	ng/L	Levels (ug/L)	大型装置							
Benzo(k)fluoranthene	0.7	10	N/A								
Benzo(a)pyrene	0.2	10	N/A								
Indeno(1,2,3-cd)Pyrene	0.4	10	N/A								
Dibenz(a,h)Anthracene	0.3	10	N/A								
benzo(g,h,i)perylene	N/A	10	N/A								

Notes:

B - Applies to organic data only. Present in the corresponding method blank.
J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.
E - Applies to Gas Chromatography/Mass Spectroscopy data only.
Indicate compound above or below linear range of instrument.

MCL - Maximum Contaminant Level (U.S. Drinking Water Standards)

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a) Blank Space - Value is below the Method Detection Limit(MDL). N/A - Not Applicable

DANNELLY ANG - Montgomery, Alabama (*)

Client Sample ID				TRAVEL BLANK	TRAVEL BLANK	TRAVEL BLANK	TRAVEL BLANK TRAVEL BLANK TRAVEL BLANK TRAVELBLK2-26 TRIP BLANK TRV-02-2-28	TRIP BLANK	TRV-02-2-28
Lab Sample Number				28998018	29046006	29017016	28909009	28891019	28934015
Matrix				WATER	WATER	WATER	WATER	WATER	WATER
Sample Date							2/26/91		2/28/91
大きない まいかんない ないない かんしゅう かんしゅ かんしゅん かんしゃ かんしゅん かんしゃ かんしゃ かんしゃ かんしゃ かんしゃ かんしゃ かんしゃ かんし	MCL	CRDL	Action						
Semivolatile Compounds	ng/L	ug/L	Levels (ug/L)			東京学芸書 第二十二十二十二十二十二十二十二十二十二十二十二十二十二十二十二十二十二十二十	表 100mm 100		
Di-n-butylphthalate	N/A	10	N/A						
N-Nitrosodiphenylamine(1)	N/A	10	N/A						
Naphthalene	N/A	10	N/A						
bis(2-Ethylhexyl)phthalate	N/A	10	3						
Phenol	N/A	10	20,000						
bis(2-Chloroethyl)Ether	N/A	10	0.03						
2-Chlorophenol	N/A	10	200						
1,3-Dichlorobenzene	009	01	N/A						
1,4-Dichlorobenzene	750	10	N/A						
Benzyl Alcohol	N/A	10	N/A			·			

B - Applies to organic data only. Present in the corresponding method blank.

J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.

E - Applies to Gas Chromatography/Mass Spectroscopy data only.

MCL - Maximum Contaminant Level (U.S. Drinking Water Standards) Indicate compound above or below linear range of instrument.

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a)

Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

DANNELLY ANG - Montgomery, Alabama (a)

Client Sample ID				TRAVEL BLANK	TRAVEL BLANK	TRAVEL BLANK TRAVEL BLANK TRAVEL BLANK TRAVELBLK2-26 TRIP BLANK TRV-02-2-28	TRAVELBLK2-26	TRIP BLANK	TRV-02-2-28
Lab Sample Number				28998018	29046006	29017016	28909009	28891019	28934015
Matrix				WATER	WATER	WATER	WATER	WATER	WATER
Sample Date							2/26/91		2/28/91
	MCL	CRDL	Action	新疆 教育 医对抗性毒素		· 在 · · · · · · · · · · · · · · · · · ·			
Semivolatile Compounds	ng/L	ng/L	Levels (ug/L)	()	があれば、数を表を認め				
1,2-Dichlorobenzene	009	10	N/A						
2-Methylphenol	N/A	10	N/A						
bis(2-Chloroisopropyl)Ether	N/A	10	N/A						
4-Methylphenol	N/A	10	N/A						
N-Nitroso-di-n-propylamine	N/A	10	0.005						
Hexachloroethane	N/A	10	20						
Nitrobenzene	N/A	10	20						
Isophorone	N/A	10	06						
2-Nitrophenol	N/A	10	N/A						
2,4-Dimethylphenol	N/A	50	N/A						

- B Applies to organic data only. Present in the corresponding method blank.
- J Applies to organic data only. Value detected is greater than zero but less than the CRDL.
 - E Applies to Gas Chromatography/Mass Spectroscopy data only.
 - Indicate compound above or below linear range of instrument.
- MCL Maximum Contaminant Level (U.S. Drinking Water Standards)
 - CRDL Contract Required Detection Limit
- Action Levels proposed in Appendix A of 40CFR254.521(a) Blank Space Value is below the Method Detection Limit(MDL).
 - - N/A Not Applicable
- (a) Data reported in ug/L.

DANNELLY ANG - Montgomery, Alabama (*)

Client Sample ID				TRAVEL BLANK	TRAVEL BLANK	TRAVEL BLANK	TRAVEL BLANK TRAVEL BLANK TRAVEL BLANK TRAVELBLK2-26 TRIP BLANK TRV-02-2-28	TRIP BLANK	TRV-02-2-28
Lab Sample Number				28998018	29046006	29017016	28909009	28891019	28934015
Matrix				WATER	WATER	WATER	WATER	WATER	WATER
Sample Date							2/26/91		2/28/91
はいません かんかん ない これ はない はない あいない	MCL	CRDL	Action						The second second
Semivolatile Compounds	ng/L	ug/L	Levels (ug/L)					1000年の大学のでは、1000年の日本の日本の日本の日本の日本の日本の日本の日本の日本の日本の日本の日本の日本の	· · · · · · · · · · · · · · · · · · ·
Benzoic Acid	N/A	10	N/A						
bis(2-Chloroethoxy)Methane	N/A	10	N/A						
2,4-Dichlorophenol	N/A	10	100						
1,2,4-Trichlorobenzene	70	10	700						
4-Chloroaniline	N/A	10	N/A						
Hexachlorobutadiene	N/A	10	4						
4-Chloro-3-methylphenol	N/A	10	N/A						
2-Methylnaphthalene	N/A	10	N/A						
Hexachlorocyclopentadiene	N/A	10	200						
2,4,6-Trichlorophenol	N/A	10	3						

B - Applies to organic data only. Present in the corresponding method blank.
J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.
E - Applies to Gas Chromatography/Mass Spectroscopy data only.
Indicate compound above or below linear range of instrument.
MCL - Maximum Contaminant Level (U.S. Drinking Water Standards)

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a)

Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable (a) - Data reported in ug/L.

QC DATA 187th Fighter Group

DANNELLY ANG - Montgomery, Alabama (*)

Client Sample ID				TRAVEL BLANK	TRAVEL BLANK	TRAVEL BLANK	TRAVEL BLANK TRAVEL BLANK TRAVEL BLANK TRAVELBLK2-26 TRIP BLANK TRV-02-2-28	TRIP BLANK	TRV-02-2-28
Lab Sample Number				28998018	29046006	29017016	28909009	28891019	28934015
Matrix				WATER	WATER	WATER	WATER	WATER	WATER
Sample Date							2/26/91		2/28/91
をできた。 こうかい かいこう かいかい かんしゅう できる できない ないしゅう かいかい かいかい かいかい かいかい かいかい かいかい かいかい かい	MCL	CRDL	Action				が、大きのでは、東京などのでは、大きのでは、大きのでは、大きのでは、大きのでは、大きのでは、大きのでは、大きのでは、大きのでは、大きのでは、大きのでは、大きのでは、大きのでは、大きのでは、大きのでは、	というない はのできる	
Semivolatile Compounds	ug/L	ng/L	Levels (ug/L)				このできるなどのでは、		
2,4,5-Trichlorophenol	N/A	50	4000						
2-Chloronaphthalene	N/A	10	N/A						
2-Nitroaniline	N/A	50	N/A						
Dimethyl Phthalate	N/A	10	N/A						
Acenaphthylene	N/A	10	N/A						
2,6-Dinitrotoluene	N/A	10	0.05						
3-Nitroaniline	N/A	50	N/A						
Acenaphthene	N/A	10	N/A						
2,4-Dintrophenol	N/A	50	70						
4-Nitrophenol	N/A	50	N/A						

Notes:

- B Applies to organic data only. Present in the corresponding method blank.
- J Applies to organic data only. Value detected is greater than zero but less than the CRDL.
 - E Applies to Gas Chromatography/Mass Spectroscopy data only.

Indicate compound above or below linear range of instrument.

MCL - Maximum Contaminant Level (U.S. Drinking Water Standards)

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a)

Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

DANNELLY ANG - Montgomery, Alabama (4)

Client Sample ID				TRAVEL BLANK	TRAVEL BLANK	TRAVEL BLANK	TRAVEL BLANK TRAVEL BLANK TRAVEL BLANK TRAVELBLK2-26 TRIP BLANK TRV-02-2-28	TRIP BLANK	TRV-02-2-28
Lab Sample Number				28998018	29046006	29017016	28909009	28891019	28934015
Matrix				WATER	WATER	WATER	WATER	WATER	WATER
Sample Date							2/26/91		2/28/91
	MCL	CRDL	Action						
Semivolatile Compounds	ug/L	ug/L	Levels (ug/L)						
Dibenzofuran	N/A	10	N/A						
2,4-Dinitrotoluene	N/A	10	N/A						
Diethylphthalate	N/A	10	30,000						
4-Chlorophenyl-phenylether	N/A	10	N/A						
Fluorene	N/A	20	N/A						
4-Nitroaniline	N/A	50	N/A						
4,6-Dinitro-2-methylphenol	N/A	50	N/A						
4-Bromophenyl-phenylether	N/A	10	N/A						
Hexachlorobenzene	1	10	N/A						
Pentachlorophenol	1	20	1000						

Notes:

B - Applies to organic data only. Present in the corresponding method blank.

J - Applies to organic data only. Value detected is greater than zero but less than the CRDL. E - Applies to Gas Chromatography/Mass Spectroscopy data only.

Indicate compound above or below linear range of instrument.

MCL - Maximum Contaminant Level (U.S. Drinking Water Standards)

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a) Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

DANNELLY ANG - Montgomery, Alabama 🗝

Client Sample ID				TRAVEL BLANK	TRAVEL BLANK	TRAVEL BLANK	TRAVEL BLANK TRAVEL BLANK TRAVEL BLANK TRAVELBLK2-26 TRIP BLANK TRV-02-2-28	TRIP BLANK	TRV-02-2-28
Lab Sample Number				28998018	29046006	29017016	28909009	28891019	28934015
Matrix				WATER	WATER	WATER	WATER	WATER	WATER
Sample Date							16/97/7		2/28/91
	MCL	CRDL	Action						ASSESSMENT OF THE PARTY OF THE
Semivolatile Compounds	ug/L	ug/L	Levels (ug/L)						
Phenanthrene	N/A	10	N/A						
Anthracene	N/A	10	N/A						
Fluoranthene	N/A	10	N/A						
Pyrene	N/A	10	N/A						
Butylbenzylphthalate	N/A	10	7,000						
3,3'-Dichlorobenzidine	N/A	20	80.0						
Benzo(a)anthracene	0.1	10	N/A						
Chrysene	0.2	10	N/A						
Di-n-octylphthalate	N/A	10	N/A						
Benzo(b)fluoranthene	0.2	10	N/A					,	

B - Applies to organic data only. Present in the corresponding method blank.

J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.

E - Applies to Gas Chromatography/Mass Spectroscopy data only. Indicate compound above or below linear range of instrument.

MCL - Maximum Contaminant Level (U.S. Drinking Water Standards)

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a) Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable (a) - Data reported in ug/L.

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DANNELLY ANG - Montgomery,
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Client Semple ID				TRAVEL BLANK	<u>TRAVEL BLANK TRAVEL BLANK TRAVEL BLANK TRAVELBLK2-26 TRIP BLANK TRV-02-2-28</u>	TRAVEL BLANK	TRAVELBLK2-26	TRIP BLANK	TRV-02-2-28
Tob Sample Number				28998018	29046006	29017016	28909009	28891019	28934015
Motrix				WATER	WATER	WATER	WATER	WATER	WATER
Sample Date							2/26/91		2/28/91
	MCL	CRDL	Action					1.1000 1000	
Semivolatile Compounds	ug/L	ug/L	Levels (ug/L)				1000年間である。	これが表表がある。	
Benzo(k)fluoranthene	0.2	10	N/A						
Benzo(a)pyrene	0.2	10	N/A						
Indeno(1,2,3-cd)Pyrene	0.4	10	N/A						
Dibenz(a,h)Anthracene	0.3	10	N/A						
benzo(g,h,i)perylene	N/A	10	N/A						

Notes:

B - Applies to organic data only. Present in the corresponding method blank.

J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.

E - Applies to Gas Chromatography/Mass Spectroscopy data only.

MCL - Maximum Contaminant Level (U.S. Drinking Water Standards)

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a)

Blank Space - Value is below the Method Detection Jimit(MDL).

N/A - Not Applicable

DANNELLY ANG - Montgomery, Alabama (*) 187th Fighter Group QC DATA

Client Sample ID				EQUIPMENT	EQUIPMENT ERB05-3-6 DUP ERB05-3-6 MS ERB-01-2-26 ERB-02-2-27 ERB-03-2-28 ERB-04-3-5 ERB-05-3-6	ERB05-3-6 MS	ERB-01-2-26	ERB-02-2-27	ERB-03-2-28	ERB-04-3-5	ERB-05-3-6
Lab Sample Number				17894020	29017P15	29017M15	28909010	28922004	28934014 28998017 29017015	28998017	29017015
Matrix				WATER	WATER	WATER	WATER	WATER	WATER	WATER	WATER
Sample Date					3/6/91	3/6/91	2/26/91	2/27/91	2/28/91	3/5/91	3/6/91
から、 一大大学 () 「 () 「 () 「 () 「 () 「 () 「 () 「 () 「 () 」 「 () 「 () 「 () 」 「 () 」 「 () 「 () 」 「 () 「 () 」 「 ()	MCL	CRDL	Action								
Semivolatile Compounds	ug/L	ng/L	Levels (ug/L)							建筑建筑	~ 法法法院
Di-n-butylphthalate	N/A	10	N/A								
N-Nitrosodiphenylamine(1)	N/A	10	N/A				2 BJ	3.3			
Naphthalene	N/A	10	N/A								
bis(2-Ethylhexyl)phthalate	N/A	10	3					10			5 BJ
Phenol	N/A	10	20,000								
bis(2-Chloroethyl)Ether	N/A	10	0.03								
2-Chlorophenol	N/A	10	200								
1,3-Dichlorobenzene	009	10	N/A								
1,4-Dichlorobenzene	750	10	N/A								
Benzyl Alcohol	N/A	10	N/A								

B - Applies to organic data only. Present in the corresponding method blank. J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.

E - Applies to Gas Chromatography/Mass Spectroscopy data only.

Indicate compound above or below linear range of instrument.

MCL - Maximum Contaminant Level (U.S. Drinking Water Standards)

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a)

Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

DANNELLY ANG - Montgomery, Alabama (*) 187th Fighter Group QC DATA

Client Sample ID				EQUIPMENT	EQUIPMENT ERB05-3-6 DUP ERB05-3-6 MS ERB-01-2-26 ERB-02-2-27 ERB-03-2-28 ERB-04-3-5 ERB-05-3-6	ERB05-3-6 MS	ERB-01-2-26	ERB-02-2-27	ERB-03-2-28	ERB-04-3-5	ERB-05-3-6
Lab Sample Number				17894020	29017P15	29017M15	28909010 28922004 28934014	28922004	28934014	28998017	29017015
Matrix				WATER	WATER	WATER	WATER	WATER	WATER	WATER	WATER
Sample Date					3/6/91	3/6/91	2/26/91	2/27/91	2/28/91	3/5/91	3/6/91
のは、日本のでは	MCL	CRDL	Action								
Semivolatile Compounds	ng/L	ng/L	Levels (ug/L)					· 大大学 经营业			100 mm
1,2-Dichlorobenzene	009	10	N/A								
2-Methylphenol	N/A	10	N/A								
bis(2-Chloroisopropyl)Ether	N/A	10	N/A								
4-Methylphenol	N/A	10	N/A								
N-Nitroso-di-n-propylamine	N/A	10	0.005								
Hexachloroethane	N/A	10	20								
Nitrobenzene	N/A	10	20								
Isophorone	N/A	10	06								
2-Nitrophenol	N/A	10	N/A								
2,4-Dimethylphenol	N/A	50	N/A								

B - Applies to organic data only. Present in the corresponding method blank. I - Applies to organic data only. Value detected is greater than zero but less than the CRDL.

E - Applies to Gas Chromatography/Mass Spectroscopy data only.

MCL - Maximum Contaminant Level (U.S. Drinking Water Standards)

Indicate compound above or below linear range of instrument.

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a) Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable (a) - Data reported in ug/L.

DANNELLY ANG - Montgomery, Alabama (a)

Client Sample ID				EQUIPMENT	EQUIPMENT ERB05-3-6 DUP ERB05-3-6 MS ERB-01-2-26 ERB-02-2-27 ERB-03-2-28 ERB-04-3-5 ERB-05-3-6	ERB05-3-6 MS	ERB-01-2-26	ERB-02-2-27	ERB-03-2-28	ERB-04-3-5	ERB-05-3-6
Lab Sample Number				17894020	29017P15	29017M15	28909010	29017M15 28909010 28922004	28934014 28998017 29017015	28998017	29017015
Matrix				WATER	WATER	WATER	WATER	WATER	WATER	WATER	WATER
Sample Date					3/6/91	3/6/91	2/26/91	2/27/91	2/28/91	3/5/91	3/6/91
	MCL	CRDL	Action				100 CO 100 CO				
Semivolatile Compounds	ug/L	ng/L	Levels (ug/L)			计操纵分析作为编制					二次 医液体管的
Benzoic Acid	N/A	10	N/A								
bis(2-Chloroethoxy)Methane	N/A	10	N/A								
2,4-Dichlorophenol	N/A	10	100								
1,2,4-Trichlorobenzene	70	10	200								
4-Chloroaniline	N/A	10	N/A								
Hexachlorobutadiene	N/A	10	4								
4-Chloro-3-methylphenol	N/A	10	N/A								
2-Methylnaphthalene	N/A	10	N/A								
Hexachlorocyclopentadiene	N/A	10	200	,							
2,4,6-Trichlorophenol	N/A	10	3								

B - Applies to organic data only. Present in the corresponding method blank.
J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.
E - Applies to Gas Chromatography/Mass Spectroscopy data only.
Indicate compound above or below linear range of instrument.
MCL - Maximum Contaminant Level (U.S. Drinking Water Standards)

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a) Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable (a) - Data reported in ug/L.

DANNELLY ANG - Montgomery, Alabama (*)

Client Samule ID				FOUTPMENT	EOUIPMENT/ERB05-3-6 DUP/ERB05-3-6 MS/ERB-01-2-26/ERB-02-2-27/ERB-03-2-28/ERB-04-3-5/ERB-05-3-6	ERB05-3-6 MS	ERB-01-2-26	ERB-02-2-27	ERB-03-2-28	ERB-04-3-5	ERB-05-3-6
I ab Sample Number				17894020	29017P15	29017M15	28909010	29017M15 28909010 28922004 28934014	28934014	28998017	29017015
Matrix				WATER	WATER	WATER	WATER	WATER	WATER	WATER	WATER
Sample Date					3/6/91	3/6/91	2/26/91	2/27/91	2/28/91	3/5/91	3/5/91
がなんでいてなる情報が最後の	MCL	CRDL	Action								
Semivolatile Compounds	ug/L	ng/L	Levels (ug/L)							(A) (基本) (基本) (基本) (基本) (基本) (基本)	が生活と言葉を
2,4,5-Trichlorophenol	N/A	20	4000								
2-Chloronaphthalene	N/A	10	N/A								
2-Nitroaniline	N/A	50	N/A								
Dimethyl Phthalate	N/A	10	N/A								
Acenaphthylene	A/N	10	N/A								
2,6-Dinitrotoluene	N/A	10	0.05								
3-Nitroaniline	N/A	50	N/A								
Acenaphthene	N/A	10	N/A								
2,4-Dintrophenol	N/A	50	70								
4-Nitrophenol	N/A	50	N/A								

B - Applies to organic data only. Present in the corresponding method blank.
 J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.
 E - Applies to Gas Chromatography/Mass Spectroscopy data only.

MCL - Maximum Contaminant Level (U.S. Drinking Water Standards) Indicate compound above or below linear range of instrument.

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a) Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

DANNELLY ANG - Montgomery, Alabama 🗝

Client Sample ID				EQUIPMENT	EQUIPMENT ERB05-3-6 DUP ERB05-3-6 MS ERB-01-2-26 ERB-02-2-27 ERB-03-2-28 ERB-04-3-5 ERB-05-3-6	ERB05-3-6 MS	ERB-01-2-26	ERB-02-2-27	ERB-03-2-28	ERB-04-3-5	ERB-05-3-6
Lab Sample Number	-			17894020	29017P15	29017M15	28909010 28922004	28922004	28934014	28998017 29017015	29017015
Matrix				WATER	WATER	WATER	WATER	WATER	WATER	WATER	WATER
Sample Date					3/6/91	3/6/91	16/97/7	2/27/91	2/28/91	3/5/91	3/6/91
The second secon	MCL	CRDL	Action	建 信							Appendig of the Appendig of th
Semivolatile Compounds	ug/L	ug/L	Levels (ug/L)							Part of the Best of	
Dibenzofuran	N/A	10	N/A								
2,4-Dinitrotoluene	N/A	10	N/A								
Diethylphthalate	N/A	10	30,000								
4-Chlorophenyl-phenylether	N/A	10	N/A								
Fluorene	N/A	50	N/A								
4-Nitroaniline	N/A	50	N/A								
4,6-Dinitro-2-methylphenol	N/A	50	N/A								
4-Bromophenyl-phenylether	N/A	10	N/A								
Hexachlorobenzene	1	10	N/A								
Pentachlorophenol	1	50	1000								
						The second secon					

- B Applies to organic data only. Present in the corresponding method blank.
- J Applies to organic data only. Value detected is greater than zero but less than the CRDL.
 E Applies to Gas Chromatography/Mass Spectroscopy data only.
 Indicate compound above or below linear range of instrument.
 MCL Maximum Contaminant Level (U.S. Drinking Water Standards)

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a) Blank Space - Value is below the Method Detection Limit(MDL). N/A - Not Applicable (a) - Data reported in ug/L.

DANNELLY ANG - Montgomery, Alabama (*)

Client Sample ID				EOUIPMENT	EQUIPMENT ERB05-3-6 DUP ERB05-3-6 MS ERB-01-2-26 ERB-02-2-27 ERB-03-2-28 ERB-04-3-5 ERB-05-3-6	ERB05-3-6 MS	ERB-01-2-26	ERB-02-2-27	ERB-03-2-28	ERB-04-3-5	ERB-05-3-6
Lab Sample Number				17894020	29017P15		28909010	28922004	29017M15 28909010 28922004 28934014 28998017 29017015	28998017	29017015
Matrix				WATER	WATER	1	WATER	WATER	WATER	WATER	WATER
Sample Date					3/6/91	3/6/91	2/26/91	2/27/91	2/28/91	3/5/91	3/6/91
ないのではないでは、 できないのできない。	MCL	CRDL	Action								
Semivolatile Compounds	ug/L	ug/L	Levels (ug/L)								
Phenanthrene	N/A	10	N/A								
Anthracene	N/A	10	N/A								
Fluoranthene	N/A	10	N/A								
Pyrene	N/A	10	N/A								
Butylbenzylphthalate	N/A	10	7,000								
3,3'-Dichlorobenzidine	N/A	20	0.08								
Benzo(a)anthracene	0.1	10	N/A								
Chrysene	0.2	10	N/A								
Di-n-octylphthalate	N/A	10	N/A								
Benzo(b)fluoranthene	0.7	10	N/A								

B - Applies to organic data only. Present in the corresponding method blank.
J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.
E - Applies to Gas Chromatography/Mass Spectroscopy data only.
Indicate compound above or below linear range of instrument.

MCL - Maximum Contaminant Level (U.S. Drinking Water Standards)

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a)

Blank Space - Value is below the Method Detection Limit(MDL). N/A - Not Applicable (a) - Data reported in ug/L.

QC DATA

187th Fighter Group

Client Sample ID Lab Sample Number Matrix Sample Date Semivolatile Compounds	MCL	CRDL	DANNE Action Levels (ug/L)	DANNELLY ANG - Montgomery, Alabama EQUIPMENT ERB05-3-6 DUP ERB05-3-6 MS ERB-01-2-26 ERB-02-2-27 ERB-03-2-28 ERB-04-3-17894020 17894020 29017P15 29017M15 28909010 28922004 289334014 28998017 Action WATER WATER WATER WATER WATER WATER WATER Action Action 3/6/91 3/6/91 3/6/91 3/5/91 3/5/91	ERB05-3-6 MS 29017M15 WATER 3/6/91	ma (a) ERB-01-2-26 28909010 WATER 2/26/91	ma (a) ERB-01-2-26 ERB-02-2-27 28909010 28922004 WATER WATER 2/26/91 2/27/91	ERB-03-2-28 28934014 WATER 2/28/91	ERB-04-3- 28998017 WATER 3/5/91
Benzo(K)Huorantnene	7.0	10	N/A						
Benzo(a)pyrene	0.2	10	N/A						

3-5|ERB-05-3-

WATER 2901701

3/6/91

A A A

의의의

0.3 N/A

ndeno(1,2,3-cd)Pyrene Dibenz(a,h)Anthracene ocnzo(g,h,i)perylene B - Applies to organic data only. Present in the corresponding method blank.

J - Applies to organic data only. Value detected is greater than zero but less than the CRDL.

E - Applies to Gas Chromatography/Mass Spectroscopy data only.

Indicate compound above or below linear range of instrument. MCL - Maximum Contaminant Level (U.S. Drinking Water Standards)

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a) Blank Space - Value is below the Method Detection Limit(MDL).

N/A - Not Applicable

QC DATA 187th Fighter Group

DANNELLY ANG - Montgomery, Alabama (*)

Client Sample ID				EOUIPMENT	EOUIPMENT ERB05-3-6 DUP ERB05-3-6 MS ERB-01-2-26 ERB-02-2-27 ERB-03-2-28 ERB-04-3-5 ERB-05-3-6	ERB05-3-6 MS	ERB-01-2-26	ERB-02-2-27	ERB-03-2-28	ERB-04-3-5	ERB-05-3-6
Lab Sample Number				17894020	29017P15	29017M15	28909010	28922004	28934014	28998017 29017015	29017015
Matrix				WATER	WATER (c)	WATER (d)	WATER	WATER	WATER	WATER	WATER
Sample Date					3/6/91	3/6/91	2/26/91	2/27/91	2/28/91	3/5/91	3/6/91
	MCL	CRDL	Action								
Inorganics	ug/L	ng/L	Level (ug/L)								京 美麗 建设
Antimony	9	09	10		N/C	93.1					
Arsenic	20	10	50		D/N	6.66					
Barium	2,000	200	1000		N/C	103.4		3.7 B			
Cadmium	5	S	10		J/N	86.8					
Chromium	100	10	50		N/C	100.6					
Copper	1,300	25	200		D/N	0.76	2.4 B	72.1			
Lead	15	3	50		N/C	92.0		21.9			
Mercury	2	0.2	2		N/C	109.0					
Nickel	100	40	700		N/C	98.6					
Selenium	50	\$	10		N/C	9.68					
Silver	N/A	01	50		N/C	92.8					
Thallium	2	10	3		N/C	100.2					
Zinc	N/A	20	2000		J/N	95.8	5.0 B	50.1			2.6 B

B - Value detected is less than the CRDL but greater than or equal to the MDL. MCL - Maximum Contaminant Level (U.S. Drinking Water Standards)

CRDL - Contract Required Detection Limit
Action Levels proposed in Appendix A of 40CFR254.521(a)
Blank Space - Value is below the Method Detection Limit(MDL).
N/C - No Change in the RPD

(c) - Relative Percent Difference(RPD) for duplicate samples.

(d) - Percent Recovery(%R) for Matrix Spike(MS) samples.

(a) - Data reported in ug/L.

QC DATA 187th Fighter Group

DANNELLY ANG - Montgomery, Alabama (*)

Client Sample ID				ERB-06-3-7	ERB-07-3-8	ER-1 F	⁷ B-01-3-4	FB-02-3-7	FB-03-3-7	TRAVEL BLANK	ERB-06-3-7 ERB-07-3-8 ER-1 FB-01-3-4 FB-02-3-7 FB-03-3-7 TRAVEL BLANK TRAVEL BLANK
Lab Sample Number				29041001	29046005 28891020 28973001 29032002 29032001	28891020	28973001	29032002	29032001	29032003	28973002
Matrix				WATER	WATER	WATER WATER	WATER	WATER WATER	WATER	WATER	WATER
Sample Date				3/7/91	3/8/91		3/4/91	3/7/91	3/7/91		
Comment of the Commen	MCL	CRDL	Action							7	
Inorganics	ug/L	ng/L	Level (ug/L)								一下 中国 中國 医二次 医二次
Antimony	9	09	10								
Arsenic	90	10	50								
Barium	2,000	200	1000					35.2 B	٠		
Cadmium	5	5	10								
Chromium	100	10	20								
Copper	1,300	25	200				51.2	22.9 B			
Lead	15	3	50				1.5 B	2.1 B	1.3 B		
Mercury	2	0.2	2								
Nickel	100	40	700								
Selenium	20	5	10								
Silver	N/A	10	20								
Thallium	2	10	3								
Zinc	N/A	20	2000			,	4.2 B	3.1 B	3.0B		

B - Value detected is less than the CRDL but greater than or equal to the MDL. MCL - Maximum Contaminant Level (U.S. Drinking Water Standards)

CRDL - Contract Required Detection Limit
Action Levels proposed in Appendix A of 40CFR254.521(a)
Blank Space - Value is below the Method Detection Limit(MDL).

N/C - No Change in the RPD

(c) - Relative Percent Difference(RPD) for duplicate samples. (d) - Percent Recovery(%R) for Matrix Spike(MS) samples. (a) - Data reported in ug/L.

DANNELLY ANG - Montgomery, Alabama (a)

Client Sample ID				TRAVEL BLANK	TRAVEL BLANK	TRAVEL BLANK	TRAVEL BLANK TRAVEL BLANK TRAVEL BLANK TRAVELBLK2-26 TRIP BLANK TRV-02-2-28	TRIP BLANK	FRV-02-2-28
Lab Sample Number				28998018	29046006	29017016	28909009	28891019	28934015
Matrix				WATER	WATER	WATER	WATER	WATER	WATER
Sample Date							2/26/91		2/28/91
	MCL	CRDL	Action						
Inorganics	ng/L	ug/L	Level (ug/L)	建筑工程设施工程					
Antimony	9	09	10						
Arsenic	20	10	20	-					
Barium	2,000	200	1000						
Cadmium	5	5	10						
Chromium	100	10	20						
Copper	1,300	25	200						
Lead	15	3	50						
Mercury	2	0.2	2						
Nickel	100	40	700						
Selenium	20	5	10						
Silver	N/A	10	50						
Thallium	2	10	3						
Zinc	N/A	20	2000						

B - Value detected is less than the CRDL but greater than or equal to the MDL. MCL - Maximum Contaminant Level (U.S. Drinking Water Standards)

CRDL - Contract Required Detection Limit

Action Levels proposed in Appendix A of 40CFR254.521(a) Blank Space - Value is below the Method Detection Limit(MDL).

N/C - No Change in the RPD (c) - Relative Percent Difference(RPD) for duplicate samples. (d) - Percent Recovery(%R) for Matrix Spike(MS) samples. (a) - Data reported in ug/L.

66-O



BORING NUMBER B1PS

SHEET

OF 1

1

PROJECT	Dannelly ANG			LOCATION	Background	
	206.2'(Gr.), 206.04'	(Toc)	DRILLING CONTRACTOR	Kilman Bros	s./Stone Mt., GA	
DRILLING ME	ETHOD AND EQUIPMENT	3¼" ID	HSA, CME 75		26/91 (1120hrs)	

WATER	LEVELS				START 2/26/91 (1020 NISH 2/26/9)	
Š.F.	5	SAMPLE		STANDARD	SOIL DESCRIPTION	COMMENTS
DEPTH BELOW SURFACE (FT)	INTERVAL	NUMBER AND TYPE	RECOVERY (FT)	PENETRATION TEST RESULTS 6"-6"-6" (N)	SOIL NAME, USCS GROUP SYMBOL, COLOR, MOISTURE CONTENT, RELATIVE DENSITY OR CONSISTENCY, SOIL STRUCTURE, MINERALOGY	DEPTH OF CASING, DRILLING RATE, DRILLING FLUID LOSS, TESTS AND INSTRUMENTATION
- 200	1.5	1S	0.8	5-5-7-7 (12)	Fat Clay (CH) - Reddish brown-orange, moist, stiff, upper 0.2' -	Chem. Analysis (1S)
_					<u>-</u>	-
				5-6-6	Giller (GT MT) Makking house appropri	·
	5.0	2S	1.5	(12)	Silty Clay (CL-ML) Mottled brown-orange—w/tan, moist, stiff—	
5 —						
-					- 	
-						(35)
-	10	3S	1.5	8-11-14 (25)	Silty Clay (CL-ML) Mottled brown-tan, - moist, v. stiff	Chem. Analysis (3S) -
10						_
-		:				_
-					· · · · · · · · · · · · · · · · · · ·	
-					Silty Clay (CL-ML) Similar to 3S, lower_	-
-	15	4 S	1.5	6-7-15 (22)	0.5' is <u>Chalk-</u> Lt. grey to Lt. brown, silty	
15						
-						
-						
-				22-45-50/4	Chalk - Olive green to brown with grey,_	Chem. Analysis (5S)
-	20	FC	, ,	(95)	moist, hard, micaceous, silty	Dup. Sample on BTEX Dup. Sample on TPH
20_	20	5S	1.5	(95)	: -	Dup. Dumple on 11
-					Boring Terminated @ 20' BGS Installed 5' 1.25" pvc screen and 15' 1.25" pvc casing. Added 7' graded sandpack, 2' bentonite seal, and grouted annulus to surface (0'-11').	Boring terminated @ 1120 hrs
-					Flush-grade piezometer completed with	-
-					steel cover and locking cap.	<u> </u>
-	-					-
					-	_
						_
	1		:		-	_
	<u> </u>	<u> </u>		<u> </u>		REV 11/89 FORM D1586



PROJECT NUMBER	
MGM27526.SI.FK	

BORING NUMBER

B2PS

SHEET

OF <u>1</u>

PROJECT	Dannelly A	ANG			LOCATION	Background		
ELEVATION	206.4'(Gr.)	206.23	(Toc)	DRILLING CONTRACTOR	Kilman Bro	s./Stone Mt. GA		
	ETHOD AND E	,	2111	ID HSA, CME 75				
					EEbral 2/	26 (01 (12105-0)	D C14-1-	

WATER	LEVELS	·			START 2/26/91 (1255hrs) INISH 2/26/9	1(1310hrs\OGGERB. Carlisle
ŠF.	,	SAMPLE		STANDARD	SOIL DESCRIPTION	COMMENTS
DEPTH BELOW SURFACE (FT)	NTERVAL	NUMBER AND TYPE	RECOVERY (FT)	STANDARD PENETRATION TEST RESULTS	OR CONSISTENCY, SOIL STRUCTURE,	DEPTH OF CASING, DRILLING RATE, DRILLING FLUID LOSS, TESTS AND INSTRUMENTATION
DEF	INT	ANC	REC (FT)	6"-6"-6" (N)	MINERALOGY	
				87	-	Boring B2PS is nexted with B1PS, therefore soil was analyzed in B1PS.
5_	5	18	1.5		Silty Clay (CL-ML)-Brown-tan mottled, moist, v. stiff	-
_					-	-
_					-	_
				9-15-19	Silty-Clay (CL-ML)-Similar to 1S	-
10_	10	2S	1.5	(30)	——————————————————————————————————————	-
			•		Boring terminated @ 10' bgs Installed 5 ft of 1.25" pvc screens and 5 ft of 1.25" pvc casing. Added 7' graded sandpack, 2' bentonite seal, and grouted remaining annulus to surface (0'-1'). Flush-grade piezometer completed with steel cover and locking cap.	Boring terminated @ 1310 hrs
-			,		-	_
L	L	L				



BORING NUMBER B3PS

SHEET

1

OF 1

PROJECT Dannelly ANG LOCATION Dackground	
ELEVATION 208.5'(Gr.), 208.31(Toc) DRILLING CONTRACTOR Kilman Bros./Stone Mt. GA	
DRILLING METHOD AND EQUIPMENT HSA, 3½" ID, CME 75	

WATER	LEVELS	·			2-27-91 2-2 START (1330 hrs) FINISH (144	7-91 0_hrs)LOGGER <u>B. Carlisle</u>
Š∈.		SAMPLE	=	STANDARD	SOIL DESCRIPTION	COMMENTS
DEPTH BELOW SURFACE (FT)	INTERVAL	NUMBER AND TYPE	RECOVERY (FT)	PENETRATION TEST RESULTS	SOIL NAME, USCS GROUP SYMBOL, COLOR, MOISTURE CONTENT, RELATIVE DENSITY OR CONSISTENCY, SOIL STRUCTURE,	DEPTH OF CASING, DRILLING RATE, DRILLING FLUID LOSS, TESTS AND INSTRUMENTATION
DEP	INTE	NUN AND	REC (FT)	6"-6"-6" (N)	MINERALOGY	
				14-6-4	Asphalt/Sand with Gravel - Road and fill material -	Asphalt noted in sample-Not submitted for analysis
2	2	18	0.5	(10)		
						•
	-			3-4-5	Poorly Graded Sand w/Gravel (Fill matl.)	Sample not submitted for analysis_
5	5	2S	0.8	(9)	Brown-black, dry, loose, gravel approx. 4mm.	due to possible asphalt content in fill material
_	-				· -	-
			:		_	_
					=	
				5-6-10	Silty Clay (CL-ML) - Lt. Brown-lt. grey, moist, v. stiff, micaceous,	Chemical analysis (3S)
10_	10	3S	1.5	(16)	slightly chalky (area the red), contains calcareous gravel in thin	·
					lenses	_
					_	
						_
		•		10-19-17	Silty Clay (CL-ML) Similar to 3S hard	Chemical analysis (4S)
15	15	4 S	1.5	(36)		
				•		_
				11-12-15	Silty Clay (CL-ML) Similar to 3S except	Boring terminated @ 30' BGS
20	20	5S	1.5	(27)	lower 0.2 contains chalk, greenish- grey, micaceous	Installed 10' of 1.25" pvc screen- and 20' of 1.25" pvc casing.
					•	Added 12.5' sandpack, 2' ben- tonite seal, and grouted annulus
						to surface (0'-15.5').
						Flush grade piezometer completed - with steel cover and locking cap.
			•	10-10-16	Silty Clay (CL-ML) Mottled tan-lt.	
25	25	6S	1.5	(26)	grey, moist, v. stiff, lower 0.5' is greenish-grey Chalk	
					_	:
				16-30-50/4	Chalk - Olive-green to grey, moist hard, calcareous gravel 2-4mm	Chemical Analysis (7S)
					maru, carcareous graver 2-4mm	
	- 30	7S	1.5	(80)		



BORING NUMBER

B5PS

SHEET 1

OF 1

PROJEC	т	Danne	lly AN	3	LOCATION	ekground			
					DRILLING CONTRACTOR Kilman Bros./St	cone Mt. GA			
DRILLING METHOD AND EQUIPMENT 3½" ID HSA, CME 75 2-26-91 2-26-91									
WATER	LEVELS	<u> </u>			2-26-91 2-26- START (1505) FINISH (1650	LOGGER B. Carlisle			
%F		SAMPLE		STANDARD	SOIL DESCRIPTION	COMMENTS			
DEPTH BELOW SURFACE (FT)	NTERVAL	NUMBER AND TYPE	RECOVERY (FT)	PENETRATION TEST RESULTS	SOIL NAME, USCS GROUP SYMBOL, COLOR, MOISTURE CONTENT, RELATIVE DENSITY OR CONSISTENCY, SOIL STRUCTURE,	DEPTH OF CASING, DRILLING RATE, DRILLING FLUID LOSS, TESTS AND INSTRUMENTATION			
DEP	INTE	AND	EEC	6"-6"-6" (N)	MINERALOGY				
_	2	18	0.9	5-6-65 (12)	Silty Clay (CL-ML) Reddish brown, moist, stiff	Chem. Analysis (1S)			
					· -				
-					-	-			
 5 —	5 ·	2S	1.5	5-3-4 (7)	Silty Clay with Sand (CL-ML) Mottled _ reddish brown-tan, moist, firm				
5 —									
-					-				
-					· -	-			
_					-	-			
-	10	38	1.2	3 -4- 6 (10)	Fat Clay with Sand (CH) Mottled - reddish brown, moist, stiff	Chem. Analysis (3S)			
10 —					, —	· 			
-					-	-			
-						-			
_									
				3-4-4	Fat Clay with Sand (CH) Similar to 3S				
15	15	4S	1.5	(8)	lower 0.5' becomes Chalky, 1t. grey-lt. brown, ferruginous, micaceous				
15 —									
-						-			
-					-	-			
4					·				
				6-19-19	Silty Clay (CL-ML) Mottled lt. brown-	_			
20	20	5S	1.5	(38)	lt. grey, moist, hard, chalky, micaceous				
20	•				<u> </u>	-			
-					-	· · · · · · · · · · · · · · · · · · ·			
-					-	· -			
_			4		-	-			
_				12-19-30	Silty Clay (CL-ML) Lt. brown-lt. grey,	Chem. Analysis (6S)			
25	25	6S	1.2	(49)	mottled, moist, hard, chalky, lower 0.2' is <u>Chalk</u> - olive-green to grey, moist, hard	Boring terminated at 30' bgs Installed 5' of 1.25" pvc screen			
-					_	and 25' of 1.25" pvc casing. Added 7' graded sandpack, 2'			
_					_	bentonite seal, and grouted			
_					·	annulus to surface (0-21').			
-	30	7S	1.5	12-21-50 (71)	Chalk - Olive green-grey, moist, hard, micaceous	Flush-grade piezometer completed with steel cover and locking cap.			



BORING NUMBER

B6MW (Monitoring Well) HEET

OF <u>1</u>

PROJECT	Dannelly ANG		LOCATION Background						
	198.4'(Gr.), 198.23'(T	OC) DRILL	ING CONTE	RACTOR Ki	lman Bros	s./Stone Mt. G	:A		.sle
	THOD AND EQUIPMENT	31" ID HSA,	CME 75						
WATER LEVE			START	2-27-91 (1005)	FINISH _	² -27-31	LOGGER _	B. Carlisle	

	LEVELS			27-91 1040) LOGGER <u>B. Carlisle</u>				
30] :	SAMPLE		STANDARD	SOIL DESCRIPTION	COMMENTS		
DEPTH BELOW SURFACE (FT)	NTERVAL	NUMBER AND TYPE	RECOVERY (FT)	PENETRATION TEST RESULTS 6"-6"-6"	SOIL NAME, USCS GROUP SYMBOL, COLOR, MOISTURE CONTENT, RELATIVE DENSITY OR CONSISTENCY, SOIL STRUCTURE, MINERALOGY	DEPTH OF CASING, DRILLING RATE, DRILLING FLUID LOSS, TESTS AND INSTRUMENTATION		
SCE	N	NA DN	EF)	(N)	MINERALOGY			
		·				Boring B6MW is nested with B5PS, therefore soil was analyzed - only in B5PS.		
5	5	15	1.5	5-6-8 (14)	Silty Clay (CL-ML)-Brown-tan w/dark organic silt, moist, stiff	_		
3 -								
					Fat Clay with Sand (CH) - Mottled			
10 _	10	2S	1.5	3-4-5 (9)	orange-brown to tan, moist, stiff, slightly chalky			
						- -		
				7-8-12	Silty Clay (CL-ML)-Brown w/lt. grey chalky laminae, moist v. stiff	- -		
15 _	15	3\$	1.5	(20)				
	-							
	20	4S	1.5	13-33-50/5 (83)	Chalk - Greenish-grey, moist, hard, silty, micaceous	Boring terminated @ 1040 hrs.		
20_								
			. .		Boring terminated @ 20' bgs Installed 10' of 2" pvc screen and 10' of 2" pvc casing. Added 12.5' graded sandpack, 2' bentonite seal, and grouted annulus to surface (0'-5.5').	_ 		
					Flush-grade monitoring well completed with steel cover, locking cap, and brass ID plate.	_		
						_		



BORING NUMBER

P1BS

SHEET 1

OF 1

PROJEC	PROJECTDannelly ANG		G	LOCATION POL (Site 1)					
ELEVATION <u>202. 6' (Gr.)</u>					DRILLING CONTRACTOR Kilman Bros./Stone Mt. GA				
DRILLING METHOD AND EQUIPMENT HSA, 3½" ID, CME 75									
WATER LEVELS					START (0800 hrs) FINISH (0855 hrs) LOGGER B. Carlisle				
ð£	SAMPLE STANDARD PENETRATION			STANDARD PENETRATION	SOIL DESCRIPTION COMMENTS				
DEPTH BELOW SURFACE (FT)	NTERVAL	NUMBER AND TYPE	RECOVERY (FT)	TEST RESULTS 6"-6"-6"	SOIL NAME, USCS GROUP SYMBOL, COLOR, MOISTURE CONTENT, RELATIVE DENSITY OR CONSISTENCY, SOIL STRUCTURE, MINERALOGY DEPTH OF CASING, DRILLING RATE, DRILLING FLUID LOSS, TESTS AND INSTRUMENTATION				
SU	Z	AN	ᆱ	(N) Not					
2		18		Sampled	Not Sampled-Limestone backfill HNU headspace @ borehole = 120ppm -				
-		2S	0.9	6-6-8 (14)	Silty Clay (CL-ML)-Mottled lt. grey-tan, Chem. Analysis (2S) moist, stiff, micaceous, slightly chalky-				
4 -				5-6-9	Silty Clay (CL-ML)-Similar to 2S Borehole headspace w/HNU = 10-15ppm				
_ 6		3S	1.2	(15)	— — — — — — — — — — — — — — — — — — —				
1 7				7-10-15	Silty Clay (CL-ML)-Similar to 2S except very stiff, less than 5% calcareous				
8		4S	1.5	(25)	gravel, clay is ferruginous				
					Silty Clay (CL-ML)-Tan with light grey, Chem. Analysis (5S) moist, v. stiff, silty TPH & BTEX Dup taken				
10_		58	1.5	(27)	_				
		6S	1.5	7-12-15 (27)	Silty Clay (CL-ML)-Similar to 5S				
12	$\overline{}$				Silty Clay (CL-ML)-Similar to 5S Chem. Analysis (7S)				
14		7S -	1.5		Lower 0.5 is Chalk-Greenish-grey-olive,- moist, hard, silty				
	Y				-				
· —					Boring terminated @ 14' bgs- Annulus grouted to surface				
-					-				
_					_Soil HNU Headspace (ppm)				
					_BS = 220ppm				
					4S = 130ppm 6S = 1ppm				
					7S = 5ppm -				
			_						
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BORING NUMBER

P2BS

SHEET

1

OF 1

			22 - 211	· · · · · · · · · · · · · · · · · · ·	POI	(Site 1)	
PROJEC	т	Danne	IIY AN	G	DRILLING CONTRACTOR Kilman Bros./Stone Mt. GA		
ELEVAT	ION	(01.8"	(6r.)	HS/	A, 3½" ID, CME 75	no ne on	
				MENI		91)LOGGER <u>B. Carlisle</u>	
		SAMPLE		STANDARD	SOIL DESCRIPTION	COMMENTS	
DEPTH BELOW SURFACE (FT)	INTERVAL	NUMBER AND TYPE	RECOVERY (FT)	PENETRATION TEST RESULTS 6"-6"-6" (N)	SOIL NAME, USCS GROUP SYMBOL, COLOR, MOISTURE CONTENT, RELATIVE DENSITY OR CONSISTENCY, SOIL STRUCTURE, MINERALOGY	DEPTH OF CASING, DRILLING RATE, DRILLING FLUID LOSS, TESTS AND INSTRUMENTATION	
				Not Sampled	Not Sampled - Limestone backfill -	-	
2 _		18		5-7-8	Silty Clay (ML-CL) - Mottled tan to	Chem. Analysis (2S)	
4_		2S	1.2		greenish-grey, moist, stiff	- -	
		3\$	1.5	4-3-4 (7)	Silty Clay (ML-CL) - Similar to 2S except firm	·	
6_				14-14-14	Silty Clay (ML-CL) - Mottled tan- lt. grey, moist, v. stiff	- -	
8_		4S	1.5	(28) 8-14-14	Silty Clay (ML-CL) - Similar to 4S	Chem. Analysis (5S)	
10		58	1.5	(28)			
		6S	1.5	4-4-9 (13)	Silty Clay (ML-CL) - Similar to 4S except stiff	-	
12_				9-12-16	Silty Clay (ML-CL) - Similar to 4S except hard, lower 0.5' is Chalk -	Chem. Analysis (7S)	
14		7\$	1.5	(38)	greenish-grey, moist	-	
_					Boring terminated at 14 ft. bgs- annulus grouted to surface	-	
					· · · · · · · · · · · · -	Soil HNU Headspace (ppm)	
					-	2S = 250 3S = 160 4S = 140	
-						4S = 140 6S = 10	
	-				-		
-							
					_		
-	-				_	-	



BORING NUMBER

P3BS

SHEET 1

OF 1

PROJEC	ст	nnelly	ANG		LOCATION POL (Site 1)			
ELEVAT	10N <u>2</u>	01.2'	(Gr.)		DRILLING CONTRACTOR Kilman Bros./Sto	one Mt. GA		
DRILLIN	IG METI	HOD AN	D EQUIF	PMENT HS	A, 3¼" ID CME 75			
WATER	LEVELS	·			2-28-91 2-28-91 START <u>(1450)</u> FINISH <u>(1545</u>) LOGGER B. Carlisle		
WO.		SAMPLE		STANDARD PENETRATION	SOIL DESCRIPTION	COMMENTS		
H BEL ACE (F	, AL	ER TYPE	VERY	TEST RESULTS	SOIL NAME, USCS GROUP SYMBOL, COLOR, MOISTURE CONTENT, RELATIVE DENSITY	DEPTH OF CASING, DRILLING RATE, DRILLING FLUID LOSS,		
DEPTH BELOW SURFACE (FT)	INTERVAL	NUMBER AND TYPE	RECOVERY (FT)	6"-6"-6" (N)	OR CONSISTENCY, SOIL STRUCTURE, MINERALOGY	TESTS AND INSTRUMENTATION		
_				Not Sampled	Not sampled-Limestone backfill -	-		
2 _		15			_	·		
				4-5-6	Silty Clay (ML-CL)-Mottled greenish- brown to lt. grey, moist, stiff,	Chem. Analysis (2S)		
4		2S	1.2	(11)	contains white chalky silt			
		38			Silty Clay (ML-CL) - Similar to 2S -	3S sampled with hand auger to check for utility obstruction		
6_		- 33		14 12 15	-			
_		40	, ,	14-13-17	Silty Clay (ML-CL)-Similar to 2S except hard -	Chem. Analysis (4S)		
8 _		4 S	1.5	(30)	-	_		
_		5S	1.5	5-6-8 (14)	Silty Clay (ML-CL) - Greenish-brown to tan, moist, stiff	-		
10_		25	1.5	4-5-8				
_		6S	1.5	(13)	Silty Clay (ML-CL) - Greenish brown-tan with lt. grey, moist, stiff	_		
12 _		05	<u> </u>	21-40-50/2	Chalk - Creenish grey-alive	Cham Analysis (7C)		
		78	1.0	(90)	Chalk - Greenish grey-olive, moist hard, silty -	Chem. Analysis (7S)		
14 _		75	1.0	(90)	-	·		
_					Boring terminated @ 14' bgs- Annulus grouted to surface			
_								
_					-	Soil HNU Headspace (ppm)		
_					· -	4S = 9		
_					-	5S = 3 6S = 30		
					·	7S = 11		
-					-	-		
_					-	· . -		
_					_	-		
-					-	_		
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PROJECT	NU	MB	ER
MCM275	26	СT	ГV

BORING NUMBER P4BS

SHEET

OF,

PROJECTDannelly ANG			lly ANG	3		OL (Site 1)
		202. 2	.' (Gr.)	DRILLING CONTRACTORKilman Bros./S	Stone Mt. GA
ELEVATION <u>202. 2' (Gr.)</u> DRILLING METHOD AND EQUIPMENT HSA,					, 3¼" ID, CME 75	-5-91
WATER LEVELS					3-5-91 3- START(1225 hrs) FINISH(13	-5-91 300-hrs) LOGGER B. Carlisle
ðf.		SAMPLE		STANDARD PENETRATION	SOIL DESCRIPTION	COMMENTS
DEPTH BELOW SURFACE (FT)	INTERVAL	NUMBER AND TYPE	RECOVERY (FT)	TEST RESULTS	SOIL NAME, USCS GROUP SYMBOL, COLOR, MOISTURE CONTENT, RELATIVE DENSITY OR CONSISTENCY, SOIL STRUCTURE,	DEPTH OF CASING, DRILLING RATE, DRILLING FLUID LOSS, TESTS AND INSTRUMENTATION
SURI	IN H	AND	REC (FT)	6"-6"-6" (N)	MINERALOGY	
				Not	Not Sampled-Limestone backfill	
2		15		Sampled		
				4-10-12	Silty Clay (CL-ML)-Mottled tan-lt. grey, moist, v. stiff, chalky -	Cham. Analysis (2S)
4	. \	2S	1.0	(22)		
_				3-3-3	Silty Clay (CL-ML)-Tan to brown w/lt. grey calcareous silt, moist, firm —	·
6		3S	1.5	(6) ⁻		-
_				6-8-12	Silty Clay (CL-ML)-Similar to 3S except v. stiff	Chem. Analysis (4S)
8		4S	1.2	(20)	<u>-</u>	
				4-4-6	Silty Clay (CL-ML)-Similar to 3S except stiff	
10	\	5 _, S	1.5	(10)		_
_				5-5-8	Silty Clay (CL-ML)-Similar to 3S except stiff, contain slight amount -	
12	\ \	6S	1.5	(13)	roots	·
_				8-15-37	Silty Clay (CL-ML)-Similar to 3S- Lower 1.0' is <u>Chalk</u> -Greenish-grey to _	Chem. Analysis (7S)
14		7S ´	1.5	(52)	olive, moist, hard, silty	
_					·	<u> </u>
-					Boring terminated @ 14' bgs- Annulus grouted to surface	
-		7.			-	
.					-	Soil HNU Headspace (ppm)
] _					_	3S = 130
						3S = 130 5S = 55 6S = 60
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PROJECT	N	UMB	EF
MGM2752	6	ST	FΚ

BORING NUMBER P5BS

SHEET 1 OF1

						(0)		
PROJECT Dannelly ANG					LOCATION POL (Site 1)			
ELEVATION <u>201. 6' (Gr.)</u> DRILLING METHOD AND EQUIPMENT HSA					DRILLING CONTRACTOR Kilman Bros./Sto	one Mt. GA		
DRILLIN WATER I				MENT HSA	2-28-91 2-28-	-91 O-hrs) LOGGER B. Carlisle		
≥	;	SAMPLE	<u> </u>	STANDARD	SOIL DESCRIPTION	COMMENTS		
DEPTH BELOW SURFACE (FT)	NTERVAL	NUMBER AND TYPE	RECOVERY (FT)	PENETRATION TEST RESULTS 6"-6"-6"	SOIL NAME, USCS GROUP SYMBOL, COLOR, MOISTURE CONTENT, RELATIVE DENSITY OR CONSISTENCY, SOIL STRUCTURE, MINERALOGY	DEPTH OF CASING, DRILLING RATE, DRILLING FLUID LOSS, TESTS AND INSTRUMENTATION		
- 200	Z	ZZ IS	 Er	(N) Not Sampled	Not sampled-Limestone backfill -	Soil descriptions above the Chalk (7S) are of POL backfill material		
2 _	$\overline{}$				No recovery			
		2S	NR		-			
4 _ 				1-1-2	Well Graded Sand with Gravel (SW) (Backfill material) tan-brown, wet,	Chem. Analysis (3S)		
6		3S	0.9	(3)	very loose			
- 8 -		4 S	1.2	(2)	Well Graded Sand with Gravel and Clay (SW-SC) - Similar to 3S escept less than 15% Clay	Backfill grain size analysis		
- -		5S	1.2	2-2-2	Well Graded Sand wih Gravel and Clay (SW-SC) - Similar to 4S	Chem. Analysis (5S)		
10		6S	1.5	(10)	Well Graded Sand with Gravel and Clay (SW-SC) - Similar to 4S except lower 0.2' is greenish-grey Chalk	-		
12 - -		7 S	1.5	8-25-30 (55)	Chalk - olive-green to lt. grey, moist hard, silty	Chem. Analysis (7S)		
14 _ - -					Boring terminated @ 14' bgs- Installed 10' of 2" pvc screen to temporarily keep the annulus open. After obtaining a water sample the screen was pulled and the remaining annulus was grouted to surface.	Could not collect Shelby tube at 7S due to refusal (hardness) *Soil HNU Hadspace (ppm) 6S = 60		
_					-	Split spoon would not retain enough sample to perform headspace analysis in the saturated backfill.		
			•		-			
					_	- -		
 					- -			
-					-			



BORING NUMBER

P6BS

SHEET 1

OF 1

	···					
PROJEC			lly AN		LOCATION PO	or (Site 1)
ELEVAT	_				DRILLING CONTRACTORKilman Bros./St	one Mt. GA
					1, 3½" ID, CME 75 3-5-91 3-	5-91
WATER	LEVELS	·			START (1005 hrs)- FINISH (111	5 hrs) LOGGER B. Carlisle
ŞΕ		SAMPLE	≣	STANDARD	SOIL DESCRIPTION	COMMENTS
DEPTH BELOW SURFACE (FT)	INTERVAL	MBER O TYPE	RECOVERY (FT)	PENETRATION TEST RESULTS 6"-6"-6"	SOIL NAME, USCS GROUP SYMBOL, COLOR, MOISTURE CONTENT, RELATIVE DENSITY OR CONSISTENCY, SOIL STRUCTURE,	DEPTH OF CASING, DRILLING RATE, DRILLING FLUID LOSS, TESTS AND INSTRUMENTATION
SCH	N T	AND AND	H. (F)	(N)	MINERALOGY	
				Not	Asphalt Pavement and backfill	
2		18		Sampled	-	
-				5-5-8	Silty Clay (CL-ML)-Mottled lt. grey-	
4		2S	1.5	(13)	tan, moist, stiff, silty	-
	7			5-7-9	Silty Clay (CL-ML)-Similar to above except v. stiff, sample contained	Intended to collect sample (3S) for analysis but the tubes
6		3S	0.8	(16)	asphalt cavings from road	contained asphalt which was carry- down from the road above.
				6-8-10	Silty Clay (CL-ML)-Similar to 2S except v. stiff	
8		4 S	1.0	(18)	v. still	Tri Sample may contain asphare
				11-14-13	Silty Clay (CL-ML)-Similar to 2S except v. stiff	Sample Analysis (5S)
10_		5S	1.2	(27)	2	
				12-10-15	Silty Clay (CL-ML)-Similar to 2S except v. stiff	· · · · · · · · · · · · · · · · · · ·
12		6S	1.5	(25)	_	_
12 -				7-11-14	Silty Clay (CL-ML)-Similar to 2S except v. stiff-Lower 0.3' is Chalk-Greenish-	Sample Analysis (7S)
14		7S -	1.2	(25)	grey-olive, silty	
					·	
					Boring terminated @ 14 ft. bgs-	
					Annulus grouted to surface	Ī
						Soil HNU Headspace (ppm)
						1S = 5
1					-	2S = 3
	ļ					5S = 3 6S = 1
-					-	<u> </u>
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BORING NUMBER

P7BS

SHEET 1

OF 1

					when the contract of the contr	
PROJEC			nnelly		LOCATION PO	OL (Site 1)
ELEVAT	$100 \frac{2}{3}$	202.4	l' (Gr.		DRILLING CONTRACTOR Kilman Bros./S	Stone Mt. GA
DRILLIN	IG MET	HOD AN	ID EQUII	PMENT HSA	, 3½" ID, CME 75 3-5-91 3-5	5-91
WATER	LEVELS	S			START (1405 hrs) FINISH (144	15 hrs) LOGGER B. Carlisle
ME(E)		SAMPL		STANDARD PENETRATION	SOIL DESCRIPTION	COMMENTS
DEPTH BEL	INTERVAL	NUMBER AND TYPE	RECOVERY (FT)	TEST RESULTS	SOIL NAME, USCS GROUP SYMBOL, COLOR, MOISTURE CONTENT, RELATIVE DENSITY	DEPTH OF CASING, DRILLING RATE, DRILLING FLUID LOSS,
DEPT	INTE	AND	E)	6"-6"-6" (N)	OR CONSISTENCY, SOIL STRUCTURE, MINERALOGY	TESTS AND INSTRUMENTATION
-		18		Not Sampled	Not Sampled-Limestone	Soil descriptions above Sample 6S are of the POL backfill material
2_	/		<u> </u>	1-1-0	Well Graded Sand with Gravel (SW) (Backfill)-wet, loose	Could not perform sample analysis due to low recovery
4_		2\$	0.2	(1)	Wall Grand G. J. Co.	
-		3S	0.2	(2)	Well Graded Sand with Gravel (SW) (Backfill)-wet, loose	Could not perform sample analysis due to low recovery
6_		33	0.2	2-1-1	Well Graded Sand with Gravel and Clay	Sample Analysis (4S)
8		4S	0.7	(2)	(SW-SC)-(Backfill)-wet, loose _	-
	Ζ,			2-1-1	Well Graded Sand with Gravel and Clay (SW-SC)-(Backfill)-wet, loose	Sample Analysis (5S)
10_		5S	0.8	(2)		
-	\	6S	0.8	6-12-50/4 (72)	Well Graded Sand with Gravel and Clay (SW-SC)-(Backfill)-wet, loose	Sample Analysis (6S)
12 _		•			Boring terminated @ 12' bgs- Hit refusal-Driller believes the auger - has encountered a concrete slab that anchors the USTs.	Could not collect Shelby Tube beneath backfill due to refusal on potential concrete slab.
					Boring terminated at 12 ft. bgs- Installed 10 ft. of 2" pvc screen to temporarily keep the annulus open. After obtaining a water sample the screen was pulled and the remaining annulus was grouted to surface.	Split spoon would not retain enough sample to perform HNU headspace.
					-	-
					-	
-			;		-	
-					-	-
						<u>-</u> -
					-	-
1					-	-



BORING NUMBER P8BS

SHEET 1

OF 1

						L (Site 1)
PROJECT					LOOATION	
ELEVAT				HSA	DRILLING CONTRACTOR Kilman Bros./S	tone Mt. GA
DRILLIN WATER			D EQUIF	PMENT HSA,	3-5-91 3-5 START(1500 hrs)_ FINISH(163	-91 5_hrs) LOGGER <u>B. Carlisle</u>
					SOIL DESCRIPTION	COMMENTS
DEPTH BELOW SURFACE (FT)	INTERVAL	NUMBER AND TYPE	RECOVERY (FT)	STANDARD PENETRATION TEST RESULTS 6"-6"-6"	SOIL NAME, USCS GROUP SYMBOL, COLOR, MOISTURE CONTENT, RELATIVE DENSITY OR CONSISTENCY, SOIL STRUCTURE, MINERALOGY	DEPTH OF CASING, DRILLING RATE, DRILLING FLUID LOSS, TESTS AND INSTRUMENTATION
<u> </u>	Ľ	ž¥	R.	(N) Not	Not Sampled-Limestone backfill	
-		1S		Sampled	NOC Sampled-Dimestone Dackilli	
2 -		2S	1.5	4-7-8 (15)	Sandy Fat Clay (CH)-Mottled brown-red, moist, stiff, lower 1.0' is Silty Clay - (CL-ML)-tan-grey, moist, stiff, chalky	Sample Analysis (2S)
6		3S	1.2	5-7-9 (16)	Silty Clay (CL-ML)-Similar to the lower 1.0 in 2S	
-		4S	1.0	7-8-9 (17)	Silty Clay-tan, moist, very stiff, silty	Sample Analysis (4S)
8 _		5S	1.5	4-7- 8 (15)	Silty Clay-tan -lt. grey, moist, stiff, silty, micaceous	
10_		6S	2.0	Shelby Tube	Silty Clay (CL-ML)-Similar to 5S as noted from open ends of Shelby Tube	SHELBY TUBE SAMPLE (6S)
14		7S -	1.0	12-22-34 (56)	Silty (CL-ML)-Similar to 5S- Lower 0.3' is Chalk Greenish-grey-olive, moist, hard, silty	Sample Analysis (7S)
-	·				Boring terminated @ 14' bgs- Annulus grouted to surface _	Soil HNU Headspace (ppm)
-				-	- -	3S = 1 ppm 4S = 1 ppm 5S = 0 ppm
-						- -
-						-
-			•		- -	-
_						_
-					- -	



BORING NUMBER 01BS

SHEET 1

OF 1

						——————————————————————————————————————
PROJEC	т	Dan	nelly	ANG	LOCATION OW	S (Site 2)
ELEVAT	ION	98.9	(Gr.)	DRILLING CONTRACTOR Kilman Bros./S	tone Mt. GA
					A, 3¼" ID, CME 75	
					3-6-91 3 START _(1000_hrs) FINISH(10	-6-91 45-hrs)_ LOGGER <u>B. Carlisle</u>
ν (T:	,	SAMPLE		STANDARD PENETRATION	SOIL DESCRIPTION	COMMENTS
DEPTH BELOW SURFACE (FT)	INTERVAL	NUMBER AND TYPE	RECOVERY (FT)	TEST RESULTS 6"-6"-6" (N)	SOIL NAME, USCS GROUP SYMBOL, COLOR, MOISTURE CONTENT, RELATIVE DENSITY OR CONSISTENCY, SOIL STRUCTURE, MINERALOGY	DEPTH OF CASING, DRILLING RATE, DRILLING FLUID LOSS, TESTS AND INSTRUMENTATION
2		18	0.8	3 - 3-3 (6)	Sandy Silty Clay (CL-ML)-Brown, moist loose, topsoil composition, organic, material with roots	Chem. Analysis (1S)
4		2S	1.5	2-2-2 (4)	Silty Clay with Sand (CL-ML)-Brown-tan with lt. grey silt, moist, soft	. -
6		3S	1.2	2-6-8	Silty Clay (CL-ML)-Mottled tan-lt. grey, moist, stiff, contains less than 5% calcareous sand	
8 _		4 S	1.5	2-4-5	Silty Clay (CL-ML)-Similar to 3S except mottled brown to red with lt. grey silt_	
10_		5S	1.5	4-5-7 (12)	Silty Clay (CL-ML)-Similar to 3S except contains roots, clay becomes chalky	- -
12		6S	1.2	(10)	Silty Clay (CL-ML)-Similar to 3S except slightly sandy (less than 10%)	-
14		7S -	1.5	8-11-10 (21)	Silty Clay with Sand (CL-ML)-Greenish- grey to brown, moist, glauconitic, sand_ less than 10%	Solvent odor noted in sample -
16 _		8S	1.5	6-11-27 (38)	Silty Clay (CL-ML)-Similar to 7S except hard - Lower 0.5' is Chalk-Greenish-grey to olive, moist, hard, silty	Chem. Analysis (8S)
-				•	Boring terminated @ 16' bgs- Annulus grouted to surface	·
-				, ,	-	Soil HNU Headspace (ppm) 2S = 1 ppm
-					_ -	5S = 40 ppm 7S = 40 ppm
			4		- -	·
_		:				
					-	- -
					_	_



BORING NUMBER

02BS

SHEET 1

OF 1

PROJECTDannelly ANG		;	LOCATION OWS (Site 2)			
ELEVATION 198.7' (Gr.)			(Gr.))	DRILLING CONTRACTOR Kilman Bros./St	one Mt. GA
DRILLING METHOD AND EQUIPMENT				MENT	HSA, 3¼" ID	(=91
WATER LEVELS					2-28-91 2-28 START (0820 hrs) FINISH (093	-91 0 hrs) LOGGER B. Carlisle
ðf;		SAMPLE		STANDARD PENETRATION	SOIL DESCRIPTION	COMMENTS
DEPTH BELOW SURFACE (FT)	NTERVAL	NUMBER AND TYPE	RECOVERY (FT)	TEST RESULTS	SOIL NAME, USCS GROUP SYMBOL, COLOR, MOISTURE CONTENT, RELATIVE DENSITY OR CONSISTENCY, SOIL STRUCTURE,	DEPTH OF CASING, DRILLING RATE, DRILLING FLUID LOSS, TESTS AND INSTRUMENTATION
SUI	Σ	AN	HE (FE)	(N)	MINERALOGY	
_		15	0.8	4-5-4 (9)	Silty Clay with Sand (CL-ML) - Tan-lt. grey, moist, stiff, contain slight organic material with roots	Chem. Analysis (IS)
2_				4-4-5	Silty Clay (CL-ML) - Tan-lt. grey, moist, stiff	:
4	. \	25	1.5	(9)	_	
		36		5-5-5	Silty Clay (CL-ML) - Greenish grey-tan, moist, stiff, slightly organic	-
6_		3S	1.5'		_	
-		4S	1.5	2 - 4-8 (12)	Silty Clay (CL-ML) - Mottled brown- tan-grey, moist, stiff, ferruginous, slightly calcareous (sand)	<u>-</u>
8_		<u> </u>		2-5-6	Silty Clay (CL-ML) - Similar to 4S	1
10		5S	0.8		-	- -
10	7			5-4-7	Silty Clay (CL-ML): - Similar to 4S	HNV scan over split spoon sample shows 5-10 ppm
12		6S	1.5	(11)	_	5110#3 3 10 ppm
				8-11-14	Silty Clay (CL-ML) - Similar to 4S	Chem. Analysis (7S) HNV scan 0-2 ppm
14		7S	1.2		Cilty Clay (CI_MI) a Similar to AS	
		8S	1.5'	12-16-32 (48)	Silty Clay (CL-ML) - Similar to 4S - lower 0.5 is Chalk-ltmed. grey, moist, hard, silty	
16_	7			16-30-50/4	Chalk - Slightly weathered tan-grey in	Chem. Analysis (9S)
18 _		9S	1.5'	(80)	upper 0.5' - lower 1.0' is greenish grey, moist, hard	· · · · · · · · · · · · · · · · · · ·
		:			Boring terminated @ 18ft bgs annulus grouted to surface	Soil HNV Headspace (ppm)
		i				2S = 0.4 3S = 0.8
					_	5S = 0.5 6S = 20
	1	i			·	8S = 22
		.]				9S = 38 -
						- -
	1				_	
		,]				·
	1					_
. 7						-
-		. [. [-1	_
-	-		.	ĺ	-	



ROJECT NUMBER	BORING NUMBE
MCM27526 ST PK	Oabc

SHEET

OF <u>1</u>

(100) IIIS)	GER <u>B. Carlisle</u> COMMENTS
NATER LEVELS START 3-6-91 FINISH 3-6-91 LOGO	
VALER LEVELS START LOGG	
	COMMENTS
SAMPLE STANDARD SOIL DESCRIPTION PENETRATION	h i
SOIL NAME, USCS GROUP SYMBOL, COLOR, MOISTURE CONTENT, RELATIVE DENSITY OR CONSISTENCY, SOIL STRUCTURE, TESTS AND I	ASING, DRILLING RATE, UID LOSS, INSTRUMENTATION
Compared to the control of the c	1- (10)
2-4-6 IS 1.0 (10) Sandy, Silty Clay (CL-ML)-Mottled brown-Chem. Analysi red, moist, stiff, organic material and roots present	IS (IS) _
2-2-2 Silty Clay (CL-ML)-Mottled brown-red,	
2S 1.5 (4) moist, soft	-
3-6-8 Silty Clay (CL-ML)-Mottled brown-tan w/ It. grey silt (calcareous), moist, stiff	•
6 Silty Clay (CL-ML)-Similar to 3S except Chem. Analysis	Ls (4S)
8 1.2 (21)	
Silty Clay (CL-ML)-Similar to 4S as noted from open ends of Shelby Tube 5S 2.0 Similar to 4S as noted from open ends of Shelby Tube	(5S) – –
8-12-23 Silty Clay (CL-ML)-Similar to 3S except hard, lower 0.2' is chalky, lt. grey to brownish-green (weathered)	-
22-35-50/5 Silty Clay (CL-ML)-Similar to lower 0.2 Chem. Analyst in 6S. Lower 1.0' is Chalk-Greenish grey to olive, moist, hard, silty	Ls (7S) - -
14	poreville Chalk
8S 3.0 CORE	
17	-
Boring terminated @ 17ft. bgs-	·
Annulus grouted to surface	-
- Soil HNU Head	ispace (ppm) -
- 2S = 5ppm 3S = 10ppm - 6S = 60ppm	
	_
-	<u></u>
4 1 -	· -
	-
	-



BORING NUMBER

04BS

SHEET

OF 1

PROJECT Dannelly ANG	LOCATION OWS (Site 2)
ELEVATION 199.2'(Gr.)	DRILLING CONTRACTOR Kilman Bros./Stone Mt. GA
DRILLING METHOD AND EQUIPMENT	HSA, 3½" ID, CME 75
WATED LEVELS	3-1-91 3-1-91 START (1030 brs) FINISH (1150) LOGGER B. Carlisle

WATER					3-1-91 3-1- START (1030-hrs) FINISH (115	0)LOGGER <u>B. Carlisle</u>
		SAMPLE		STANDARD	SOIL DESCRIPTION	COMMENTS
DEPTH BELOW SURFACE (FT)			RECOVERY (FT)	PENETRATION TEST RESULTS	SOIL NAME, USCS GROUP SYMBOL, COLOR, MOISTURE CONTENT. RELATIVE DENSITY	DEPTH OF CASING, DRILLING RATE, DRILLING FLUID LOSS, TESTS AND INSTRUMENTATION
DEPTI	INTERVAL	NUMBER AND TYPE	RECC (FT)	6"-6"-6" (N)	OR CONSISTENCY, SOIL STRUCTURE, MINERALOGY	
				3-5-6	Sandy Organic Topsoil Upper 0.4'- Silty Clay (ML-CL) - Lower 0.8'-	Chem. Analysis (1S)
2		18	1.2	(11)	mottled tan-lt grey, moist, stiff	;
				5-4-5	Silty Clay (ML-CL) - Similar to lower -	_
4	. /	2S	1.5	(9)	_	·
				4-4-7	Silty Clay (ML-CL) - Similar to lower 0.8' of 1S except slight calcareous sand content	_
6_		3 S	1.5	(11) 4-8-11	Silty Clay (ML-CL) - Similar to lower	Chem. Analysis (4S)
-		4 S	1.0	(19)	0.8' of 1S except v. stiff, chalky	-
8_				Shelby	Silty Clay (ML-CL) - Similar to lower 0.8' of 1S as noted from open ends of	Shelby Tube (5S)
10		5S	2.0	Tube	Shelby Tube	
10_	/			8-10-16	Silty Clay (ML-CL) - Tan-lt. grey, moist, v. stiff, chalky (weathered)	
12	\	6S	1.5	(26)	slight organic material noted (roots)	
				4-8-24	Silty Clay(ML-CL) - Similar to 6S except hard - No organics noted	
14		7S	1.5		lower 0.2' is <u>Chalk</u> , greenish grey, silty, hard	(0C)
-				28-35-38	Chalk w/interlayered silty clay. Chalk is olive green to grey, moist, hard	Chem. Analysis (8S) —
16		8S	1.5	(73)		_
-				·	Boring terminated @ 16' bgs- Annulus grouted to surface	-
-					-	
-					-	Soil HNU Headspace (ppm)
-						2S = 230
-					_	3S = 20 6S = 300
-					· · · · ·	7S = 150 _
-	1				-	*Note-Boring redrilled and
-					-	sampled 3/6/91 due to sample _ shipping problems
-	1					-
-	1				_	-
	1				-	
	1				_	-
	-			-	_	
L	<u> </u>	1	<u> </u>	1	1	REV 11/89 FORM D1586



BORING NUMBER 05BS

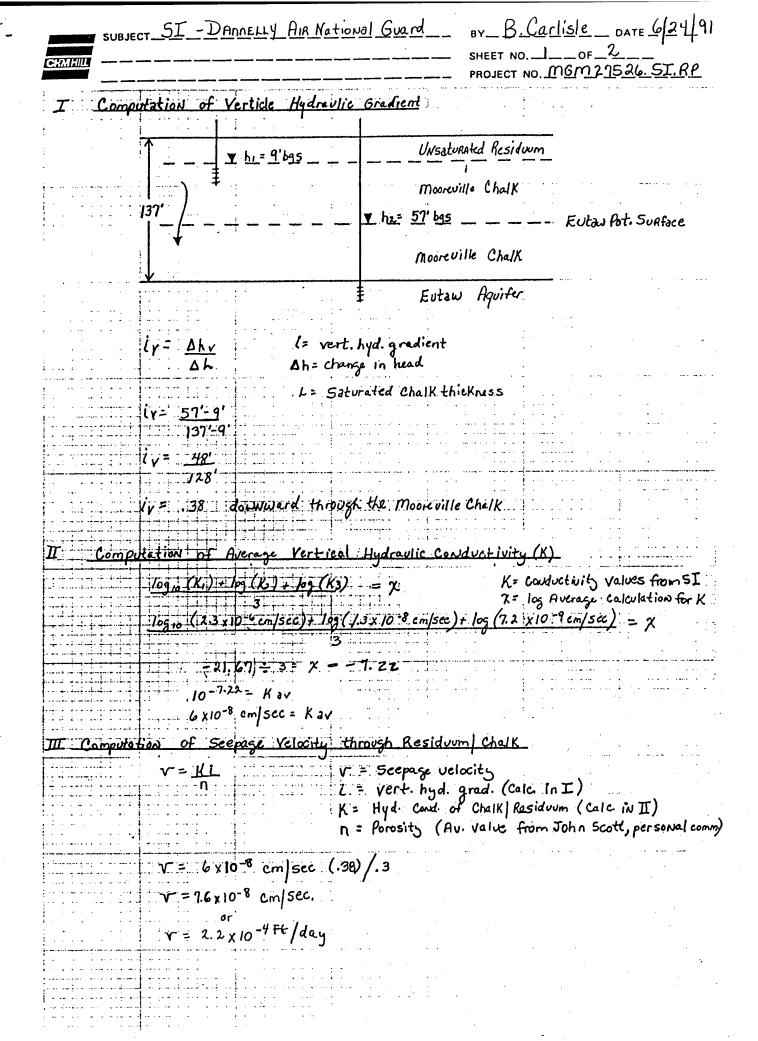
SHEET 1

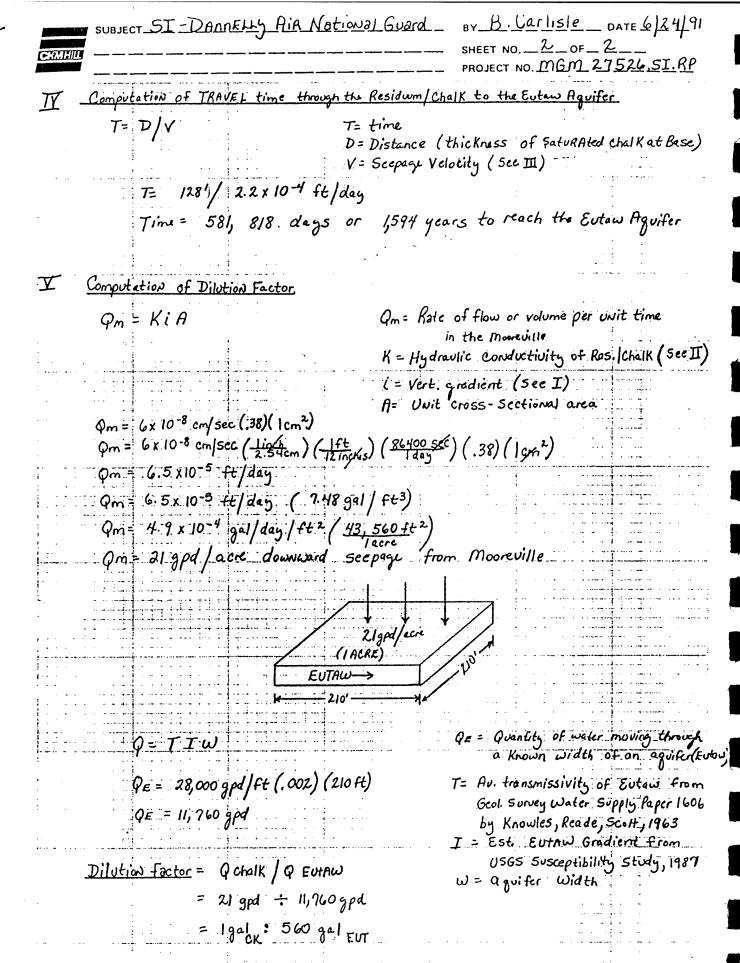
OF 1

PROJECT	Dannelly ANG	LOCATION OWS (Site 2)
ELEVATION _	199.1' (Gr.)	DRILLING CONTRACTOR Kilman Bros./Stone Mt. GA
DRILLING ME	THOD AND EQUIPMENT _	HSA, 3¼" ID, CME 75
WATED LEVE	c	3-6-91 3-6-91

				-WEINT	3-6-91 3-6 START (1310_hrs) FINISH (135	0 hrs) LOGGER B. Carlisle
		SAMPLE		STANDARD	SOIL DESCRIPTION	COMMENTS
DEPTH BELOW SURFACE (FT)	INTERVAL	NUMBER AND TYPE	RECOVERY (FT)	PENETRATION TEST RESULTS 6"-6"-6" (N)	SOIL NAME, USCS GROUP SYMBOL, COLOR, MOISTURE CONTENT, RELATIVE DENSITY OR CONSISTENCY, SOIL STRUCTURE, MINERALOGY	DEPTH OF CASING, DRILLING RATE, DRILLING FLUID LOSS, TESTS AND INSTRUMENTATION
2		1S	0.9	3-4-4	Sandy Silty Clay (CL-ML)-Brown-tan w/ lt. grey silt, moist, firm, organic w/ - roots	Chem. Analysis (1S)
4		2S	1.5	2-2-2 (4)	Silty Clay with Sand (CL-ML)-Brown-tan, wet, soft, organic material and roots - present	
6		3S	1.5	2-5-4 (9)	Silty Clay (CL-ML)-Mottled tan-lt. grey, moist, stiff, clacareous silt/sand — present	
8		4S	1.5	2-3-2 (5)	Silty Clay (CL-ML)-Similar to 3S except firm	-
		5S	1.0	4-4-8	Silty Clay (CL-ML)-Similar to 3S	Chem. Analysis (5S)
10		6S	1.5	5-6 - 10 (16)	Silty Clay (CL-ML)-Tan-lt. grey, moist, v. stiff, thin chalky laminae present _	
14		7S -	1.2	11-12-15 (27)	Silty Clay (CL-ML)-Mottled tan-brown w/lt. grey silty laminae, moist, v. stiff, lower 0.2' is Chalk-Brown-greenish grey, slightly sandy-silty	Chem. Analysis (7S)
					Boring terminated @ 14' bgs- Annulus grouted to surface _	
-					-	Soil HNU Headspace (ppm) 3S = 5
					- 	4S = 11 6S = 110
-					- -	- -
-					-	· -
						·
-					_	-
					- -	
	L		1			DEV 11/00 FORM D1500

COMPUTATION OF VERTICAL HYDRAULIC GRADIENT AND SEEPAGE VELOCITY





DATA REVIEW AND VALIDATION PROCEDURES

TO:

Margaret Corey

FROM:

Ann Castleberry

Ann West/WDC

DATE:

January 7, 1992

SUBJECT:

Data Review and Validation for Dannelly ANG

PROJECT: MGM27526.SI.OC

1. INTRODUCTION

Soil and water samples were collected as part of the Dannelly Field, Alabama Air National Guard Site Investigation (SI). The purpose of this memorandum is to summarize the criteria used and the results of the review and validation process. The data results are discussed in the main body of the SI report and are not included in this memorandum. Data validation is the technical review of a data package using criteria established in the Data Quality Objectives of the Quality Assurance Project Plan.

All the samples were submitted to and analyzed by two CH2M HILL laboratories located in Montgomery, Alabama, and Redding, California.

2. DATA PACKAGE DELIVERABLES

When samples were submitted to the laboratory, they were assigned 8-digit unique numerical sample identifiers. The first 5 digits of the laboratory sample number identify the sample batch, and the last 3 numbers indicate each unique field sample. Attachment 1 is a summary of all the field samples submitted to the two laboratories, and the corresponding laboratory numerical sample identifier and the requested analytical parameters.

As indicated in Attachment 1, samples were submitted for either HAZWRAP Level B or C QC. For this specific project, only TPH was analyzed using Level B QC; Level C data package deliverables were provided for the other analytical methods.

MEMORANDUM

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Level B data package deliverables included:

- Sample results
- Method blank results
- Initial calibration data
- Continuing calibration data
- Spiked sample results

Level C data package deliverables are summarized in Table 1.

3. LEVEL B REVIEW CRITERIA

Level B data did not undergo a formal validation process; rather, these data were reviewed for compliance with holding time, calibration, and completeness criteria as outlined below.

HOLDING TIME

Holding time is defined as the time, in days, from sample collection to sample analysis. The holding time for TPH analysis is 28 days.

CALIBRATION

Initial calibration data are used to demonstrate that the analytical instrument is performing satisfactorily and is capable of producing acceptable quantitative data at the beginning of the analytical period. Continuing calibration checks document that the instrument continues to produce acceptable data.

TPH calls for an initial 3-point calibration with a zero intercept. Samples with a TPH concentration greater than the highest calibration standard were diluted and reanalyzed. Continuing calibration checks were performed after every 10th sample.

COMPLETENESS

Completeness can have two meanings. First, it can mean that all the data package deliverables are present and reviewed. Second, completeness can be expressed as the percentage of measurements made that are judged to be valid.

Table 1 Level C Data Package Deliverables

ORGANICS--GC/MS

Form	Purpose
I	Sample results
II	Surrogate spike results
III	MS/MSD spike results
IV	Method blank data
V	GC/MS tuning data
VI	Initial calibration data
VII	Continuing calibration data
VIII	Internal standard area data
ANICSGC	
I	Sample results
II	Surrogate spike results
III	MS/MSD spike results
IV	Method blank data

METALS

VI

VII

I	Sample results
II	Initial and continuing calibration data
III	Method blank results
IV	ICP interference check sample results
V	Spike recovery data
VI	Duplicate sample results
VII	Laboratory control sample results
V.II	Standard addition results
X	Holding times

Initial calibration data

Continuing calibration data
Second column confirmation data

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4. LEVEL C REVIEW AND VALIDATION CRITERIA.

Samples that were analyzed using CLP methods were reviewed and validated using CLP guidelines. When non-CLP methods were used, the data were reviewed and validated in the same manner using laboratory-specific acceptance criteria.

ORGANIC ANALYSES

Organic data were generated using CLP methods and were reviewed and validated using the guidance document "Laboratory Data Validation Functional Guidelines for Evaluating Organics Analyses," 1988 revision. As mentioned above, when non-CLP methods were used, the data were reviewed and validated in the same manner using laboratory-specific acceptance criteria. This guidance document lists criteria for evaluating the data package form by form. The raw experimental data were summarized and presented in the appropriate form because no raw data were included in the data package.

Also included with the sample results are Tentatively Identified Compounds (TICs) results. When compounds not on the target compound list (TCL) are detected, the instrument searches its mass spectre database looking for the closest match. Because TICs are not included in the calibration process and are only identified by the mass spectre, they are considered to be tentatively identified. Additionally, quantitation of TICs is approximate because it is not known how efficiently the compound responds to the detector. Therefore, TICs are considered to be indicative of a class or type of compound at an approximate concentration, but calibration is necessary for greater definition.

Form I, Data Results

This form presents the sample results and the information necessary for calculating holding times, and is also reviewed for completeness. Holding time is defined as the time, in days, from sample collection to sample extraction/analysis. It is important to note that the holding time for extraction often is different for water and soil samples for the same analytical method. Holding times are summarized in Table 2.

Form II, Surrogate Recovery

This form summarizes the surrogate spike recovery information. Surrogate spike recoveries are used to demonstrate laboratory performance and to evaluate matrix interference. Surrogate compounds are the structural homologs of target list

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Form V, GC/MS Tuning and Mass Calibration

This form presents the tuning and mass calibration information for each GC/MS instrument used to produce data for the sample delivery group. The CLP SOW establishes tuning and performance criteria in order to ensure mass resolution, identification, and to some degree sensitivity. These criteria are not sample specific; conformance is determined using standard materials at least once during each 12-hour analytical period. Therefore, these criteria should be met in all circumstances. Bromofluorobenzene (BFB) is used for volatile analyses, and Decafluorotriphenylphosphine (DFTPP) is used for semivolatile analyses. Gas chromatographs are calibrated and not tuned, so there is no tuning or mass calibration information for GC analyses.

Form VI, Initial Calibration

This form is used to report compound recoveries from the initial calibration solutions. Initial calibration data are used to demonstrate that the analytical instrument is performing satisfactorily and is capable of producing acceptable quantitative data at the beginning of the analytical period. Initial calibration is performed in accordance with the applicable analytical method. Acceptance limits are also defined by the analytical method, when recoveries do not fall within the acceptance limits, analysis of field samples is stopped, the problem identified and resolved, and the instrument recalibrated before sample analysis begins.

Form VII, Continuing Calibration

This form is used to report continuing calibration check sample recoveries. Continuing calibration checks document that the instrument continues to produce acceptable data. Continuing calibration check samples are analyzed at a frequency required by the analytical method, and at least immediately before and after an analytical period. Acceptance limits are also defined by the analytical method. When recoveries do not fall within the acceptance limits, analysis of field samples is stopped, the problem identified and resolved, and the instrument recalibrated before sample analysis begins.

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Form VIII, Internal Standards

This form presents the internal standards peak area information. Internal standard compounds are used to ensure that instrument sensitivity and response are stable during each analytical sequence. Acceptance criteria are defined either in the CLP SOW or the specific analytical method.

INORGANIC ANALYSES

All the inorganic data were generated using CLP methods and were reviewed and validated using the guidance document "Laboratory Data Validation Functional Guidelines for Evaluating Inorganic Analyses," 1988 revision. This guidance document lists criteria for evaluating the data package form by form. The raw experimental data were summarized and presented in the appropriate form since no raw data were included in the data package.

Form I, Sample Results

The form is similar to the organic form in that it summarizes the sample results; however, a separate form (Form X) is used to calculate sample holding time.

Form II, Initial and Continuing Calibration Verification

This form is similar to the organic form in that it summarizes the calibration results and is used to evaluate initial and continuing calibration. Form IIB is used to demonstrate that the laboratory was capable of analyzing below the CRDL at the time of analysis. The laboratory analyzes a standard solution that is twice the instrument detection limit (IDL) to verify the linearity of the instrument at low detection limits.

Form III, Blanks

This form is used to report analyte concentrations detected in the initial calibration blank (ICB), continuing calibration blanks (CCB), and the preparation blanks (PB). As mentioned in the organics section, method blanks are used to monitor the presence and magnitude of contamination introduced during the analytical process.

One method blank was analyzed for every 20 samples, or one per analytical batch, whichever was more frequent. ICBs, CCBs, and PBs were analyzed at the frequency required by the CLP SOW.

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Form IV, ICP Interference Check Sample (ICS)

This form is used to report ICS results for each ICP instrument used. The ICS sample is a mixture of analytes that have a potential for interference and is performed to verify the laboratory's interelement and background correction factors.

Form VA, Pre-digestion Spike Recovery

This form is used to report results for the pre-digestion spike recovery. This spike recovery measurement is analogous to the MS/MSD in that it provides a measure of the effects of the specific sample matrix on the sample results. Additionally it provides a measure of the efficiency of the digestion process. If the pre-digestion spike recovery does not fall within the 75 to 125 percent recovery window, then a post-digestion spike is added and the sample reanalyzed.

Form VB, Post-Digestion Spike Recovery

This form is used to report post-digestion spike results when they are necessary. The acceptance limits for post-digestion spike recoveries are also 75 to 125 percent. If both the pre- and post-digestion spike reveries are outside these acceptance limits, then this is considered positive evidence of matrix interference and the data are flagged appropriately.

Form VI, Duplicates

This form is used to report duplicate laboratory results rather than field duplicate sample results. Duplicate laboratory samples differ from MS/MSD samples in that the duplicate sample is not spiked; therefore, precision must be estimated using native rather than spiked results. For this reason, laboratory duplicates were performed only on field samples and not field QC samples.

Form VII, Laboratory Control Sample (LCS) Results

This form is used to report the recovery results for the standard LCS. The LCS analysis is designed to monitor the efficiency of the digestion process. Analyte recoveries must fall with 90 to 110 percent; if not, the data are flagged.

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Form VIII, Standard Addition Results

This form is used to report the results of samples analyzed using the method of standard addition and is used only for Graphite Furnace Atomic Absorption analyses. Duplicate injections and furnace post-digestion spike recoveries are used to establish the precision and accuracy of the individual analytical determinations. Samples must be analyzed using the method of standard addition if the analyte concentration is greater than five times the CRDL. The CLP SOW requires that the results agree within + 20 percent relative standard difference or else the data is qualified.

Form IX, ICP Serial Dilution Results

ICP serially diluted samples are used to monitor whether significant physical or chemical interferences exist as a result of sample matrix effects. The sample is diluted and the results compared (diluted versus undiluted) for agreement for any analyte whose concentration is 50 times greater than the IDL.

Form X, Holding Times

This form is used to report holding times for mercury and cyanide analyses. Sample results that are not analyzed within the holding time are flagged to indicate a low bias.

QUALIFYING FLAGS

Samples that did not meet the acceptance limit criteria were qualified with a flag, single letter abbreviations that indicate a problem with the data. Although the flags originate in the data validation section, they are included in the data summary tables (in the main body of the text) so that data will not be used indiscriminately. Flags used in this text include:

- Undetected. Analyte was analyzed for but not detected above the method detection limit.
- B The analyte was detected in both the field sample and the corresponding method blank.
- <u>J</u> Estimated. The analyte was present, but the reported value may not be accurate or precise.

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It is important to note that the data summary form, Form I, may also have flags on them and the usage is the same as above with two major additions. For organic results, data that are above the method detection limit but below the contract required detection limit (CRDL) are flagged with a J. For inorganic results, if the analyte was detected above the instrument detection limit, but below the CRDL, then the result is flagged with a B.

5. RESULTS AND CONCLUSIONS

The data were reviewed and validated as indicated in the preceding sections. As each data package was reviewed, a worksheet was completed for each analysis. A blank worksheet is shown in Figure 1. These worksheets were developed to act as a checklist for the data reviewer and are included in Attachment 2 to this appendix. Any non-conformances with the data package were noted on the worksheet and then appropriate flags were assigned to the data.

Acetone and methylene chloride are used as extraction solvents; hence, they are common laboratory contaminants. When detected in a sample, the concentration reported is the actual concentration in the sample; blank subtraction (concentration in the corresponding method blank minus the concentration in the sample) was not used. Therefore all the Acetone and methylene chloride detected in these specific samples can be attributed to laboratory contamination.

Five organic compounds were detected in the subsurface soil samples collected at Site 2:

- Acetone
- Carbon disulfide
- Di-n-butylphthalate
- Methylene chloride
- bis(2-ethylhexyl)phthalate

These compounds were detected in the corresponding field and laboratory blanks; therefore, they can be attributed to field and laboratory contamination.

The data are acceptable as is and can be used in the decision-making process without further qualification.

DATA VALIDATION

CASE:SITE:	ANALYSIS:
LAB NOTES:	
·	
HOLDING TIMES:	
NUMBER OF SAMPLES: Soil	
CALIBRATION Initial:	
COEFFICIENT:	
DI ANVC	
Field:	
FIELD DUPLICATES:	
SURROGATES:	······································
MS/MSD:	
INTERNAL STANDARDS:	TUNING:
OTHER:	
SUMMARY:	
	·

Attachment 1 DATA INVENTORY TABLE

DANNELLY ANG
FIELD EFFORT
DATA INVENTORY
ATTACHMENT 1
10-Jun-91

Time: 02:07 PM

Filename: 27526COC.WK1

							$\overline{}$			$\overline{}$												0000000	-		
DATE	SUBMITTED	TO LAB	02/22/91	02/22/91	02/22/91	02/22/91	02/22/91	02/22/91	02/22/91	02/22/91	02/22/91	02/22/91	02/22/91	02/22/91	02/22/91	02/22/91	02/22/91	02/22/91	02/22/91	02/22/91	02/22/91		02/22/91	02/22/91	02/22/91
	BTEX																								
	PAH		×	×	×	×	×	×	×	×	×	×	×	×	×	×	×	×	×	×	×				
	ТРН																						×	×	×
SOLUABLE	METALS																								
TOTAL	METALS																							·	
	SV	CLP																					,		
	VOC	CLP								·															

MOM	LAB	NUMBER	17894-001	17894-002	17894-003	17894-004	17894-005	17894-006	17894-007	17894-008	17894-009	17894-010	17894-011	17894-012	17894-013	17894-014	17894-015	17894-016	17894-017	17894-018	17894-020		17898-001	17898-002	17898-003
MDM	LAB	MATRIX NUMBER	SOIL 17894-001				SOIL 17894-005				SOIL 17894-009	SOIL 17894-010			SOIL 17894-013				SOIL 17894-017	-			SOIL 17898-001	17898-002	SOIL 17898-003
MGM	QC LAB	RIX																						17898-002	_
MGM		MATRIX	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	TIOS	TIOS	SOIL	SOIL	SOIL	WATER		SOIL	SOIL 17898-002	SOIL
FIELD	8	LEVEL MATRIX	C SOIL	C SOIL	02/21/91 C SOIL	C SOIL	C SOIL	DUP 02/21/91 C SOIL	C SOIL	C SOIL	C SOIL	1 02/21/91 C SOIL	02/21/91 C SOIL	t 02/21/91 C SOLL	C SOIL	C SOIL	02/22/91 C SOIL	с ѕоп	TIOS	C SOIL	C WATER		C SOIL	02/21/91 C SOLL 17898-002	с ѕоп

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	FIELD				MGM			TOTAL	SOLUABLE				DATE
	SAMPLE	DATE	8		LAB	VOC	SV	METALS	METALS	TPH	PAH	BTEX	SUBMITTED
- 1	А	SAMPLED	LEVEL	MATRIX	NUMBER	CLP	CLP						TO LAB
	A7-SS-6-24DUP	02/21/91	ပ	SOIL	17898-004					×			02/22/91

m	B3P-S-8.5-10	02/27/91	В	SOIL	17931-001					×			02/28/91
<u> </u>	B3P-S-13.5-15	02/27/91	В	SOIL	17931-002					×			02/28/91
m	B3P-S-28.5-30	02/27/91	Д	SOIL	17931-003					×			02/28/91
A	B1P-S-0-2	02/26/91	Д	SOIL	17937-001					×			02/27/91
B	B1P-S-8.5-10	02/26/91	В	SOIL	17937-002								02/27/91
H	B1P-S-18.5-20	02/26/91	В	SOIL	17937-003								02/27/91
H	B1P-S-18.5-20DUP	02/26/91	В	SOIL	17937-004					×			02/27/91
E S	B5P-S-0-2	02/26/91	В	SOIL	17937-005					×			02/27/91
8	B5P-S-8.5-10	02/26/91	В	SOIL	17937-006					×			02/27/91
8	B5P-S-23.0-25	02/26/91	В	SOIL	17937-007					×			02/27/91
8	B5P-S-23.0-25DUP	02/26/91	В	SOIL	17937-008								02/27/91
FI	TRAVEL BLANK	02/26/91	В	SOIL	17937-009								02/27/91
띖	ERB-01-2-26	02/26/91	Д	SOIL	17937-010								02/27/91
8	02B-S-0-2	02/28/91	В	SOIL	17946-001					×			03/01/91
잗	02B-S-12-14	02/28/91	В	SOIL	17946-002					×			03/01/91
징	02B-S-18-20	02/28/91	В	SOIL	17946-003					×			03/01/91
8	P2B-S-2-4	02/28/91	Д	SOL	17946-004					×			03/01/91
21	P2B-S-8-10	02/28/91	м	SOIL	17946-005					×			03/01/91
81	P2B-S-8-10DUP	02/28/91	щ	SOIL	17946-006					×			03/01/91

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	FIELD				MGM			TOTAL	SOLUABLE				DATE
SITE	SAMPLE	DATE	8		LAB	VOC	SV	METALS	METALS	TPH	PAH	BTEX	SUBMITTED
	a	SAMPLED	LEVEL	MATRIX	NUMBER	CLP	CL						TO LAB
-	P2B-S-12-14	02/28/91	Д	SOIL	17946-007					×			03/01/91
-	P3B-S-2-4	02/28/91	щ	SOIL	17946-008					×			03/01/91
1	P3B-S-6-8	02/28/91	e Ø	SOIL	17946-009					×		·	03/01/91
1	P3B-S-12-14	02/28/91	Д	SOIL	17946-010					×			03/01/91
1	P5B-S-4-6	02/28/91	Д	SOIL	17946-011					×			03/01/91
	P5B-S-8-10	02/28/91	ф	SOIL	17946-012					×			03/01/91
1	P5B-S-12-14	02/28/91	В	SOIL	17946-013					×			03/01/91
				•									
2	04BS-0-2	03/01/91	В	SOIL	17947-001					×			03/01/91
2	04BS-6-8	03/01/91	В	SOIL	17947-002					×			03/01/91
2	04BS-14-16	03/01/91	В	SOIL	17947-003					×			03/01/91
2	04BS-14-16DUP	03/01/91	В	SOIL	17947-004		:		·	x			03/01/91
1	P1BS-2-4	03/05/91	В	SOIL	17966-001					X			03/05/91
1	P1BS-8-10	03/05/91	В	SOIL	17966-002					×			03/05/91
1	P1BS-8-10DUP	03/05/91	В	SOIL	17966-003					×			03/05/91
1	P1BS-12-14	03/05/91	В	SOIL	17966-004					×			03/05/91
-	P6BS-6-8	03/05/91	В	SOIL	17966-005					×			03/05/91
1	P6BS-8-10	03/05/91	В	SOIL	17966-006					×			03/05/91
-	P6BS-12-14	03/05/91	В	SOIL	17966-007					×			03/05/91
	P7BS-6-8	03/05/91	В	SOIL	17966-008		1			×			03/05/91
1	P7BS-8-10	03/05/91	В	SOIL	17966-009					×			03/05/91
	P7BS-12-14	03/05/91	В	SOIL	17966-010					×			03/05/91

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		1	r		ŧ ·		,		100000	,		,	r		·	1		r	r		r		,	roceres.	·
DATE	SUBMITTED	TO LAB	03/05/91	03/05/91	03/05/91	03/05/91	03/05/91	03/05/91		03/06/91	03/06/91	03/06/91	03/06/91	03/06/91	03/06/91	03/06/91	03/06/91	03/06/91	03/06/91	03/06/91	03/06/91	03/06/91	03/06/91		03/07/91
	BTEX																								
	PAH																								
	TPH		×	×	×	×	×	×		×	×	×	×	×	×	×	×	×	×	×	×	×	×		×
SOLUABLE	METALS				:																				
TOTAL	METALS																								
	SV	CLP																							
	VOC	CLP																							
MGM	[VB	NUMBER	17966-011	17966-012	17966-013	17966-014	17966-015	17966-016		17978-001	17978-002	17978-003	17978-004	17978-005	17978-006	17978-007	17978-008	17978-009	17978-010	17978-011	17978-012	17978-013	17978-014		17995-001
		MATRIX	TIOS	SOL	TIOS	SOIL	SOIL	SOIL		SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOL	SOIL	TIOS	SOIL	SOIL	SOL		SOIL
	8	LEVEL	В	В	В	В	В	Д		В	В	æ	В	В	В	В	В	В	В	В	В	В	В		Д
	DATE	SAMPLED	03/05/91	03/05/91	03/05/91	03/05/91	03/05/91	03/05/91		03/06/91	03/06/91	03/06/91	03/06/91	03/06/91	03/06/91	03/06/91	03/06/91	03/06/91	03/06/91	03/06/91	03/06/91	03/06/91	03/06/91		03/07/91
FIELD	SAMPLE	А	P4BS-2-4	P4BS-6-8	P4BS-12-14	P8BS-2-4	P8BS-6-8	P8BS-12-14		04BS-0-2	04BS-6-8	04BS-14-16	04BS-14-16DUP	01BS-0-2	01BS-10-12	01BS-14-16	05BS-0-2	05BS-8-10	05BS-12-14	03BS-0-2	03BS-6-8	03BS-12-14	03BS-12-14DUP		A1BS-0-2
	SITE		-	1	1	1	1	1		2	2	2	2	2	2	2	2	2	2	2	2	2	2 (4

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	BTEX																	×	×	×	×
	PAH																				
	TPH		×	×	×	×	×	×	×	×	×										
SOLUABLE	METALS											×					×				
TOTAL	METALS											х	×	×	×		×				
	SV	CLP																			
	voc	CLP										×	×	×	×	×	×				
MGM	LAB	NUMBER	17995-002	17995-003	17995-004	17995-005	17995-006	17995-007	18000-001	18000-002	18000-003	18301-001	18301-004	18301-006	18301008	18301-009	18311-001	28891-001	28891-002	28891003	28891-004
		MATRIX	тоѕ	SOIL	SOIL	SOIL	TIOS	TIOS	SOL	SOIL	SOL	WATER	WATER	WATER	WATER	WATER	WATER	SOL	SOIL	SOIL	TIOS
	8	LEVEL	В	В	В	В	В	В	В	щ	۰щ	Ü	ບ	υ	ပ	υ	В	ນ	၁	บ	۵
•	DATE	SAMPLED	16/1/0/60	03/07/91	03/07/91	03/07/91	03/07/91	03/07/91	03/08/91	16/80/20	03/08/91	04/10/91	04/10/91	04/10/91	04/10/91	04/10/91	04/12/91	02/22/91	02/22/91	02/22/91	02/22/91
FIELD	SAMPLE	Œ	A1BS-8-10	A1BS-16-18	A3BS-0-2	A3BS-8-10	A3BS-18-20	A3BS-18-20DUP	A2BS-4-6	A2BS-12-14	A2BS-18-20	BG-MW-6	BG-MW-6-DMS	BG-MW-6-FB	BG-MW-6-ER	TRAVEL BLANK	BG-2P	A1-SS-0-6	A1-SS-6-24	A2-SS-0-6	A2-SS-6-24
	SITE		4	4	4	7	4	. †	4	4	4	BKGD	ВКОD	BKGD	BKGD	BKGD	BKGD	4	4	4	4

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	FIELD				MGM			TOTAL	SOLUABLE				DATE
SITE	SAMPLE	DATE	8		LAB	VOC	SV	METALS	METALS	ТРН	PAH	BTEX	SUBMITTED
	А	SAMPLED	LEVEL	MATRIX	NUMBER	CLP	CLP						TO LAB
4	A3~SS-0-6	02/22/91	υ	SOIL	28891-005			,				×	04/23/91
4	A3-SS-0-6DUP	02/22/91	ပ	SOLL	28891-006							×	04/23/91
4	A3-SS-6-24	02/22/91	3	TIOS	28891-007							×	04/23/91
4	A3-SS-6-24DUP	02/22/91	C	TIOS	28891-008							×	04/23/91
4	A4-SS-0-6	02/22/91	٥	тоѕ	28891-009							×	04/23/91
4	A4-SS-6-24	02/22/91	၁	поѕ	28891-010							x	04/23/91
4	A5-SS-0-6	02/22/91	۵	TIOS	28891-011							×	04/23/91
4	A5-SS-6-24	02/22/91	၁	SOIL	28891-012							×	04/23/91
4	A6-SS-0-6	02/22/91	υ	SOIL	28891-013							×	04/23/91
4	A6-SS-6-24	02/22/91	င	SOL	28891-014							×	04/23/91
4	A7-SS-0-6	02/22/91	υ	SOIL	28891015							×	04/23/91
4	A7-SS-6-24	02/22/91	င	SOIL	28891-016							×	04/23/91
4	A8-SS-0-6	02/22/91	υ	SOIL	28891-017							×	04/23/91
4	A8-SS-6-24	02/22/91	ບ	SOIL	28891-018							×	04/23/91
8	Trip Blank	02/22/91	၁	WATER	28891-019							×	04/23/91
သွ	ER-1	02/22/91	Ü	WATER	28891-020							×	04/23/91
ВКС	B1P-S-0-2	02/26/91	ပ	SOIL	28909-001	×	×	×					02/27/91
BKGD	B1P-S-8.5-10	02/26/91	υ	SOIL	28909-002	×	×	×					02/27/91
вкар	B1P-S-18.5-20	02/26/91	ပ	SOIL	28909-003	×	X	×					02/27/91
BKGD	B1P-S-18.5-20DUP	02/26/91	υ	SOIL	28909-004	×	×	×					02/27/91
вкар	B5P-S-0-2	02/26/91	ပ	SOIL	28909-005	×	×	×					02/27/91
вкар	B5P-S-8.5-10	02/26/91	υ	SOIL	28909-006	×	×	×					02/27/91

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DATE	SUBMITTED	TO LAB	02/27/91	02/27/91	02/27/91	02/27/91	02/28/91	02/28/91	02/28/91	02/28/91		03/01/91	03/01/91	03/01/91	03/01/91	03/01/91	03/01/91	03/01/91	03/01/91	03/01/91	03/01/91	03/01/91	03/01/91	03/01/91
	BTEX														×	×	×	×	×	×	×	×	×	×
	PAH																							
	TPH												,											
SOLUABLE	METALS															-								
TOTAL	METALS		×	×		×	×	×	×	×				×	·									
	SV	CLP	×	×		×	×	×	×	×		×	×	×		,			·					
٠	VOC	CLP	×	×	×	×	×	×	×	×		×	×	×										
MGM	LAB	NUMBER	28909-007	28909-008	28909-009	28909-010	28922-001	28922-002	28922-003	28922-004		28934-001	28934-002	28934-003	28934-004	28934-005	28934-006	28934-007	28934-008	28934-009	28934-010	28934-011	28934-012	28934-013
		MATRIX	SOIL	SOIL	WATER	WATER	SOIL	SOIL	SOIL	WATER		TIOS	SOIL	SOIL	SOIL	SOIL	soir	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL
	8	LEVEL	ບ	ပ	ບ	υ	υ	υ	U	υ		ບ	υ	ပ	υ	Ü	υ	υ	υ	υ	υ	υ	υ	υ
e .	DATE	SAMPLED	16/97/20	02/26/91	02/26/91	02/26/91	16/12/70	02/27/91	02/27/91	16/12/70		02/28/91	02/28/91	02/28/91	02/28/91	02/28/91	02/28/91	02/28/91	02/28/91	02/28/91	02/28/91	02/28/91	02/28/91	02/28/91
FIELD	SAMPLE	А	BSP-S-23.0-25	B5P-S-23.0-25DUP	Travel Blank 2-26	ERB-01-2-26	B3P-S-8.5-10	B3P-S-13.5-15	B3P-S-28.5-30	ERB-02-2-27		02B-S-0-2	02B-S-12-14	02B-S-18-20	P2B-S-2-4	P2B-S-8-10	P2B-S-8-10DUP	P2B-S-12-14	P3B-S-2-4	P3B-S-6-8	P3B-S-12-14	P5B-S-4-6	P5B-S-8-10	P5B-S-12-14
	SITE		ВКО	ВКО			BKGD	BKGD	ВКС	S,		2	7	2	1		-	1		1	1	1	-	1

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DATE	SUBMITTED	TOLAB	03/01/91	03/01/91	03/05/91	03/05/91	16/50/60	03/06/91	03/06/91	03/06/91	03/06/91	03/06/91	03/06/91	03/06/91	03/06/91	03/06/91	03/06/91	03/06/91	03/06/91	03/06/91	03/06/91	
	BTEX		×	×	×			×	×	×	×	×	×	×	×	×	×	×	×	×	×	
	PAH				×																	
	TPH																					
SOLUABLE	METALS																					
TOTAL	METALS					×																
	·	CLP				×																
	VOC	CLP				×	×															
MGM	LAB	NUMBER	28934-014	28934-015	28972-001	28973-001	28973-002	28998-001	28998-002	28998-003	28998-004	28998-005	28998-006	28998-007	28998-008	28998-009	28998-010	28998-011	28998-012	28998-013	28998014	
		MATRIX	WATER	WATER	WATER	WATER	WATER	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	SOIL	
	8	LEVEL	ပ	υ	В	၁	۲	υ	c	υ	υ	υ	υ	υ	ŭ	υ	υ	၁	C	S	ນ	
	DATE	SAMPLED	02/28/91	02/28/91	03/04/91	03/04/91	03/04/91	03/05/91	03/05/91	03/05/91	16/50/50	03/05/91	03/05/91	03/05/91	03/05/91	03/05/91	03/05/91	03/05/91	03/05/91	03/05/91	03/05/91	
FIELD	SAMPLE	О	ERB-03-2-28	TRV-02-2-28	P5-BW1	FB-01-3-4	Travel Blank	P1BS-2-4	P1BS-8-10	P1BS-8-10DUP	P1BS-12-14	P6BS-6-8	P6BS-8-10	P6BS-12-14	P7BS-6-8	P7BS-8-10	P7BS-12-14	P4BS-2-4	P4BS-6-8	P4BS-12-14	P8BS-2-4	
	SITE		Ş	oc.		χ,	00	1	1	1	-	-	1	-	1	1	1	1 1	1 1	1	1	

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	FIELD				MGM			TOTAL	SOLUABLE				DATE
SITE	SAMPLE	DATE	8		EV3	voc	AS.	METALS	METALS	TPH	PAH	BTEX	SUBMITTED
	А	SAMPLED	LEVEL	MATRIX	NUMBER	CLP	CLP.						TOLAB
1	P8BS-12-14	03/05/91	၁	SOIL	28998-016							×	03/06/91
8	ERB-04-3-5	03/05/91	ນ	WATER	28998-017							×	03/06/91
8	Travel Blank	03/05/91	υ	WATER	28998-018							×	03/06/91
2.	04BS-0-2	03/06/91	၁	SOIL	29017-001	×	×						03/07/91
2	04BS-6-8	03/06/91	၁	NOIL	29017-002	×	×	×					03/07/91
2	04BS-14-16	03/06/91	၁	SOIL	29017-003	×	×						03/07/91
2	04BS-14-16DUP	03/06/91	၁	тоѕ	29017-004	×	×						03/07/91
2	01BS-0-2	03/06/91	ວ	тоѕ	29017-005	×	×						03/07/91
2	01BS-10-12	03/06/91	၁	тоѕ	29017-006	×	×						03/07/91
2	01BS-14-16	03/06/91	ລ	TIOS	29017-007	×	×	×					03/07/91
2	05BS-0-2	03/06/91	υ	SOIL	29017-008	×	×	×					03/07/91
7	05BS-8-10	16/90/60	υ	TIOS	29017-009	×	×						03/07/91
2	05BS-12-14	03/06/91	υ	SOIL	29017-010	×	×						03/07/91
7	03BS-0-2	03/06/91	ن	TIOS	29017-011	×	×						03/07/91
2	03BS-6-8	16/90/80	۵	тоѕ	29017-012	×	×						03/07/91
2	03BS-12-14	03/06/91	٥	тоѕ	29017-013	×	×	×					03/07/91
2	03BS-12-14DUP	03/06/91	၁	TIOS	29017-014	×	×	×					03/07/91
ညွ	ERB-05-3-6	03/06/91	ט	WATER	29017-015	×	×	×					03/07/91
သွ	Travel Blank	03/06/91	Ü	WATER	29017-016	×							03/07/91
oc oc	FB-03-3-7	16/1/0/80	ວ	WATER	29032-001	×	×	×					03/08/91
ပွ	FB-02-3-7	03/07/91	υ	WATER	29032-002	×	×	×					03/08/91

.

Date: Time:

> DANNELLY ANG FIELD EFFORT

Filename: 27526COC. WK1

10-Jun-91 02:07 PM

> DATA INVENTORY ATTACHMENT 1 10-Jun-91

DATE	SUBMITTED	TO LAB	03/08/91	03/08/91	03/08/91	03/08/91	03/08/91	03/08/91	03/08/91	03/08/91	03/08/91	03/08/91	03/08/91	03/08/91	03/09/91	03/09/91	03/09/91	03/09/91	03/09/91	03/09/91
	BTEX		×	×		×	x	×	×	×	×		×	×	×	×	×	×	×	×
	PAH		×	×		X	×	×	×	×	×	×	×	×	×	×	×	×		
	TPH																			
SOLUABLE	METALS																			
TOTAL	METALS																			
	SV	CLP																		
	voc	CLP			×					,										
MGM	LAB	NUMBER	29032-003	29032-004	29032-005	29032-006	29032-007	29032-008	29032-009	29032-010	29032-011	29032-012	29041-001	29041-002	29046-001	29046-002	29046-004	29046-005	29046-003	29046-006
		MATRIX	WATER	WATER	WATER	WATER	WATER	WATER	WATER	WATER	WATER	WATER	SOIL	SOIL	SOIL	SOIL	SOIL	WATER	SOIL	WATER
	8	LEVEL	υ	ပ	ပ	ပ	ပ	၁	C	၁	၁	ວ	٥	2	υ	υ	υ	Ü	ပ	ပ
	DATE	SAMPLED	03/07/91	03/07/91	03/07/91	03/07/91	03/07/91	03/07/91	03/07/91	03/07/91	03/07/91	03/07/91	03/07/91	03/07/91	03/08/91	03/08/91	03/08/91	03/08/91	03/08/91	03/08/91
FIELD	SAMPLE	А	ERB-06-3-7	P7BW	Travel Blank	A1BS-0-2	A1BS-8-10	A1BS-16-18	A3BS-0-2	A3BS-8-10	A3BS-18-20	A3BS-18-20DUP	ERB-06-3-7	P7BW	A2BS-4-6	A2BS-12-14	A2BS-18-20	ERB-07-3-8	A2BS-12-14DUP	TRAVEL BLANK
	SITE		႘	1	ညွ	4	4	4	4	4	4	4	- ي	<u>-</u> ک	4	4	4	20	4	% %

Attachment 2 DATA VALIDATION WORKSHEETS

CASE: 17894 SITE: 4 ANALYSIS: PAH
LAB NOTES: 3 le extractions were past holding times (-5, -6, -7)
HOLDING TIMES: 2/25 (extract) - 2/21 (sampling) = OK
NUMBER OF SAMPLES: Soil /8 Water / - ERB
CALIBRATION Initial: ok
Continuing: 8K
COEFFICIENT:
BLANKS Method: nothing found
Field: nothing found
V FIELD DUPLICATES: hotel dups are at least 10x smaller conc. Then notice
SURROGATES: weter ok soil - 73-115 ok terp d 14 ag. = ok
MS/MSD: good
INTERNAL STANDARDS:TUNING:
OTHER:
SUMMARY: MGM doesn't have the co-elution problem RDD has.
The M. runa concentrations were smaller than the original conc. which
may have been due to non-homogenesty of sample +/or holding time
expiration - I used cost of original sample analysis.

CASE: /7898 SITE:	4 ANALYSIS: TPH
LAB NOTES: none	
HOLDING TIMES: 3/18 an.	- 2/21 me = 34 days.
NUMBER OF SAMPLES: Soil	L Water
CALIBRATION Initial:	· · · · · · · · · · · · · · · · · · ·
Continuing: ok	·.
COEFFICIENT: 0.9998	
BLANKS Method: no h. L.	
•	
FIELD DUPLICATES:	
SURROGATES: not found	
MS/MSD:	
INTERNAL STANDARDS:	TUNING: LCS ok
OTHER:	
SUMMARY: RESULTS 1 - U	2-32 mg/kg 3-11.0 mg/kg
•	4-9.1- B1-U-

CASE: <u>/79</u>	3/ SITE:	BKGD	AN.	ALYSIS: TPH
LAB NOTES:	no problems			
	•			
HOLDING T	MES: 3/20	- 2/27	= 21 day	14
NUMBER OF	SAMPLES: So	011 <u>3</u>	Wat	er
CALIBRATIO Initial	e of icv		***	
Continu	ing: ok (C	/		
COEFFICIEN	T: 0.9998			
BLANKS Method:	nothing four	nd		
	none			
FIELD DUPI	ICATES: non	<u>د</u>		
SURROGATES	: —			
MS/MSD:	94.6 % R			
	TANDARDS:		_ TUNING:	LCS-LOR
OTHER:	on the section of the			
SUMMARY:	3 samples	-001 5	.4 mg/K	3
r	esults:	-002 5	72 "	with 10x dilution
		-003 2	8.4	

CASE: 17937 SITE: BKGD	ANALYSIS:	TPH
LAB NOTES:		* ***
HOLDING TIMES: 3/20 analysis - 2/26 sar	pled = ok	28 day
NUMBER OF SAMPLES: Soil 7		
CALIBRATION I(V-ok Initial: CCV = ok		
Continuing: ((/-e/K		
COEFFICIENT: 0.9999 ok		
BLANKS Method: no hits		
Field: norte		
FIELD DUPLICATES: # 0%D		·
SURROGATES: NA		
MS/MSD: no data		
INTERNAL STANDARDS: NA TUN	ing: LCS-, ok	
OTHER: -2 + - 3 not listed on inventory but is		Cot C.
SUMMARY: There were concentrations detect	ted in each sa	nple-
range 3.6 - 10.0 mg/kg		

CASE: 17946 SITE: 2 + 1 ANALYSIS: 7PH
LAB NOTES: no problems
HOLDING TIMES: 3/20 analysis - 2/28 sampled = 20 days
NUMBER OF SAMPLES: Soil /3 Water
CALIBRATION Initial: ICV - ok
Continuing: CCV - ok
COEFFICIENT: 0.9999
BLANKS Method: no hits
Field: none
FIELD DUPLICATES: -5 +-6 not too good = U and 49.3
SURROGATES:
MS/MSD: 98.6 70R = OR
INTERNAL STANDARDS:
OTHER:
SUMMARY: Sample results: 4 non-detects
range - 4 - 297 mg/kg
4 ora 50; 5 under
In order. 16, U, U, 189, U, #\$ 49.3, 4, 297, 196, U, 205, 33.9, 4.5

CASE: 17947 SITE: 2	ANALYSIS: TPH
LAB NOTES: no problems	· · · · · · · · · · · · · · · · · · ·
·	
HOLDING TIMES: $3/20 - 3/1$	= 19 days
HOLDING TIMES: $3/20 - 3/1$ NUMBER OF SAMPLES: Soil 4	Water
CALIBRATION CV - TK	
Continuing: <u>CCV</u> - ok	
COEFFICIENT: 0.9999 r ²	
BLANKS Method: no futo	
Field:	
FIELD DUPLICATES:	
SURROGATES:	
MS/MSD: 93.170R	
INTERNAL STANDARDS:	TUNING: LCS of
OTHER:	
SUMMARY:	
Sample -1 12.8 mg/Kg	
-2 95.1	
-3 37.8	
-4 11.3	

OLDING TIMES: 3/27 enelysis - 3/5 sempling : 22 UMBER OF SAMPLES: Soil /6 Water ALIBRATION Initial: /CV sk Continuing: CCV sk OEFFICIENT: 0.9999 LANKS Method: nothing found Field: none IELD DUPLICATES: sk URROGATES: S/MSD: 93.7 %K NTERNAL STANDARDS: TUNING: LCS	TPH
ALIBRATION Initial: 100 sk Continuing: 000 sk OEFFICIENT: 0.9999 LANKS Method: nothing found Field: none IELD DUPLICATES: sk URROGATES: S/MSD: 93.7 % K NTERNAL STANDARDS:	
ALIBRATION Initial: 100 sk Continuing: 000 sk OEFFICIENT: 0.9999 LANKS Method: nothing found Field: none IELD DUPLICATES: sk URROGATES: S/MSD: 93.7 % K NTERNAL STANDARDS:	
ALIBRATION Initial: 100 sk Continuing: 000 sk OEFFICIENT: 0.9999 LANKS Method: nothing found Field: none IELD DUPLICATES: sk URROGATES: S/MSD: 93.7 % K NTERNAL STANDARDS:	
ALIBRATION Initial: 100 sk Continuing: 000 sk OEFFICIENT: 0.9999 LANKS Method: nothing found Field: none IELD DUPLICATES: sk URROGATES: S/MSD: 93.7 % K NTERNAL STANDARDS:	days.
ALIBRATION Initial: 100 sk Continuing: 000 sk OEFFICIENT: 0.9999 LANKS Method: nothing found Field: none IELD DUPLICATES: sk URROGATES: S/MSD: 93.7 % K NTERNAL STANDARDS:	,
OEFFICIENT: 0.9999 LANKS Method: nothing found Field: none IELD DUPLICATES: ok URROGATES: S/MSD: 93.7 %K NTERNAL STANDARDS: TUNING: LCs	
LANKS Method: nothing found Field: none IELD DUPLICATES: sk URROGATES: S/MSD: 93.7 % K NTERNAL STANDARDS:	
Method:notling found Field:none IELD DUPLICATES:dk URROGATES:	
Field: IELD DUPLICATES: URROGATES: S/MSD:	
URROGATES: S/MSD: 93.7 %K NTERNAL STANDARDS:TUNING: \(\alpha \sigma \sigma \).	
S/MSD: 93.7 %K NTERNAL STANDARDS:TUNING: \(\alpha \sigma \sigma \).	
NTERNAL STANDARDS: TUNING: CS	
THER:	ه ر
	124.4
UMMARY: six samples - U; seven samples belo	w 10 mg/K
UMMARY: six samples - U; server samples selon Three big hits1 (585), -11 (538), -12 (275))

ASE: 17978 SITE: 2 ANALYSIS: TPH
AB NOTES: No Connents
OLDING TIMES: 3/28 - 3/6 = 22 days
UMBER OF SAMPLES: Soil /4 Water
ALIBRATION Initial: ICV - ok
Continuing: ccv - ok
OEFFICIENT: 0.9997 ok: r ²
ILANKS Method: nothing found Field: none
TIELD DUPLICATES: 2k
URROGATES:
15/MSD: 98 % he ok
nternal standards: Tuning : LCS = ok
THER:
SUMMARY: Sample 1 - big hit 2120 mg/kg 2 69.1
10 12.1
4,8,9 below 7
the 8 others - U

CASE: 17995 SITE: 4	analysis: TPH
LAB NOTES: No comments	
HOLDING TIMES: 3/28 - 3/07 =	21 days
NUMBER OF SAMPLES: Soil 7	
CALIBRATION Initial: 1CV 102.5 % ok	
Continuing: CCV 102.9%K -6K	
COEFFICIENT: 0.9997 12	
BLANKS Method: no lite	
Field: none	
FIELD DUPLICATES: -6 +-7 no lits	
SURROGATES:	
MS/MSD: 100.67R	
	TUNING: LCS = 100.6% - 0k
OTHER:	
SUMMARY: no lits	

CASE: 18000 SITE: 4	ANALYSIS: TPH
LAB NOTES: Method 418, 1 (MOD)	
	.
HOLDING TIMES: 3/31 (Pralypis) -	3/08 (sampling) = 23
NUMBER OF SAMPLES: Soil 3	Water
CALIBRATION Initial: 1CV - ok	
Continuing: &k	
COEFFICIENT: 8.9998 - ok	
BLANKS Method: <u>no dotects</u>	
Field:mone	
FIELD DUPLICATES: Mone	
SURROGATES:	
MS/MSD: 97.2%R-OR	
INTERNAL STANDARDS:	TUNING:
SUMMARY: no detects	
· · · · · · · · · · · · · · · · · · ·	

CASE: 18301 SITE: BK&D	ANALYSIS: VOC
LAB NOTES:	
HOLDING TIMES:	
NUMBER OF SAMPLES: Soil	Water 5 / sample
CALIBRATION Initial: •k	I FB I ER I TRIP
Continuing: ok son dates analysed	
only empdo out were not in	∨ samples
BLANKS Field Method: MeCl 3-22 Acetone 11-18 To	luene 2
method Field: MeC1 1-2 Acetone 9-14 4.1	methyl-2-Pentenone 2-2 Xylene 1
FIELD DUPLICATES:	2 Hexanone -
SURROGATES: 0k d-8 Toluese wide fluctu	ation but within 3 sd.
MS/MSD: ok	
INTERNAL STANDARDS: ok 1	CUNING: ok .
OTHER:	
SUMMARY:	
•	

18301 TOTAL + SOL CASE: <u>18300</u> SITE: <u>BKGD</u> ANALYSIS: <u>INORGANIC</u> S
LAB NOTES: Se - 1:5 dilution due to background interference
spike R out
7/25 analysis
HOLDING TIMES: 4/24 prep - 4/10 samp = 14 days
NUMBER OF SAMPLES: Soil Water 5 1 so(
CALIBRATION I ER I FB
Continuing: ok
COEFFICIENT:
Method: O.1 Hg Pb 2.9 Ba 2.2 Cr 3.6
Field: Ba 2.7 Cr 2.7 Pb 1.3 Az 10.5
FIELD DUPLICATES: similar Lab duplicates - all below TRDL
SURROGATES:
MS/MSD: Se - 17.8 % R prespike - no problem - both below RDL Se - post-dis. spike - ,81.8 70R for the soluble sample; 78.8 for total 30-
INTERNAL STANDARDS: 1CS-ok TUNING:
OTHER: LC5 - Se = 84.1 % R - 91.8% R - ok (Limits = 78-113)
SUMMARY:

CASE: 1831/ SITE: BKGD ?	ANALYSIS: VOC
AB NOTES: no matrix spike, no exce	eptions
HOLDING TIMES: 4/21 analysis -	4/12 sampled = 9 days
NUMBER OF SAMPLES: Soil	Water/
CALIBRATION Initial: MeC 24.7 % RSD	
Continuing: several over 25%) - 1	Me C1 · 28.2 - "J"
COEFFICIENT:	
BLANKS Method: MeCl - 1 Aceton 12	
Field: none	
FIELD DUPLICATES: none	
surrogates:	
MS/MSD: none	
INTERNAL STANDARDS:	TUNING:
OTHER:	
SUMMARY: no hits except the two lat	contaminanto

CASE:18311 SITE: BK6D	SOL + TOTAL ANALYSIS: INORGANIC
LAB NOTES: More	
LAB NOTES. More	
HOLDING TIMES: Hg 4/25 - 4/12 - 1	13 days
NUMBER OF SAMPLES: Soil	Water _/
CALIBRATION Initial: <u>%</u>	· .
Continuing: ok	
COEFFICIENT: LOS LUDANACE - OR	
COEFFICIENT: for furnace - ok	
BLANKS Method: mo lits	
Field: none	
FIELD DUPLICATES: none	
SURROGATES: -	
	Anna Anna Anna Anna Anna Anna Anna Anna
MS/MSD: <u>not requested</u>	
INTERNAL STANDARDS:	FUNING: LCS- all in
OTHER:	
SUMMARY: no hits for total and solub	le metals
,	
0 1 t + - 1/40 mm has/-	
Conductivity = 1640 umhos/cm	

CASE: 2889/ SITE: 4	ANALYSIS: BTEX
LAB NOTES: 1000 IS resposes caused 3 surrogate 9	Oho to be out - 3 were re-enalyted
HOLDING TIMES: 2/27 analysis - 2/2	sampled = ok
NUMBER OF SAMPLES: Soil 18	Water 2 = 1 Trip + 1 ERB
CALIBRATION Initial: K	
Continuing: ok except for TBME	E ToD but no lita
COEFFICIENT: ok for 4/1	
BLANKS Method: no hits	
Field: ms hes	
FIELD DUPLICATES: no hits	
SURROGATES: el ligh - range = 102-18	4 averaging about 140
MS/MSD: good	
INTERNAL STANDARDS: 75 = 1000	TUNING:
OTHER:	
SUMMARY: no hit	
1235 4	



BTEX SURROGATE RECOVERY

Instrument ID: GC-3600

Date		Laboratory Reference No.	EPA-602/8020 Surrogate Recovery Fluorobenzene
2-25-91		Method Blank 28891 - 1	126 132
2-25-91		28891 - 2	130
2-25-91		28891 - 3	151
2-25-91		28891 - 4	132
2-25-91		28891 - 5	142
2-25-91		28891 - 6	147
2-26-91		28891 - 7	135
2-26-91		28891 - 8	140
2-26-91		28891 - 11	160 *
2-26-91		28891 - 12	166 *
2-26-91	1	28891 - 13	169 *
2-26-91		28891 - 14	144
2-26-91		Method Blank	132
2-26-91		28891 - 15	149
2-26-91		28891 - 16	142
2-26-91		28891 - 17	148
2-26-91		28891 - 18	143
2-26-91		28891 - 8 MS	. 102
2-26-91		28891 - 8 MSD	110
2-26-91		28891 - 19	130
2-27-91		28891 - 20	129
2-27-91		Method Blank	125
2-27-91		28891 - 9	145
2-27-91		28891 - 10	140
2-27-91		28891 - 11-RE	184 *
2-28-91		28891 - 12-RE	155 *
2-28-91		28891 - 13-RE	163 *
		4 (1)	
		İ	1

EPA 602/8020 surrogate standards reported as percent recovery. NA = Not Analyzed.

Comments: * Surrogate outside control limits.

Approved By:

jas25

CASE: 28909 SITE: Background ANALYSIS: SVOC
LAB NOTES: 0-001 needed dilution
method blank contamination
3) MS/MSD from 29017-004 MUAL EXT SEMP HOLDING TIMES: 3/20 - 2/28 - 2/26 - 0K
NUMBER OF SAMPLES: Soil
CALIBRATION Initial: OK
Continuing: 70D Benzoic Acid: 999 - 4-Nitrophenel 28.2; Hexacles- +
cyclopentadine 36.8. 2, 4. Dinitalphenol 27.5, Berzo Eghid parrylene 51.8 46. Dinitro = metyl prenol 30.7 + 44.4) 1.2-Diphenyl mydragine 29.2 COEFFICIENT: NA
BLANKS Method: Water - N-nitrosodiphenylamine - give .010 a B
Field: no conc. Action: Remove B from Soil wy conc over 610
FIELD DUPLICATES: 007+008 not too similar -007 has 5 hits 008=1
SURROGATES: (Ness than 1/2 of con hit
MS/MSD:
INTERNAL STANDARDS: ok TUNING: ok
OTHER:
SUMMARY:

*

attacher page

7B SEMIVOLATILE CONTINUING CALIBRATION CHECK

Lab Name: CH2M HILL/LRD Contract: S28909

Instrument ID: 4600 Calibration date: 03/22/91 Time: 1503

Lab File ID: 91M1BN1765 Init. Calib. Date(s): 03/04/91 03/05/91

Min RRF50 for SPCC(#) = 0.050

Max %D for CCC(*) = 25.0%

COMPOUND	RRF	RRF50	%D ======
Phenol	0.804	0.894	
bis(2-Chloroethyl)Ether	0.705	0.815	-15.6
2-Chlorophenol	0.674	0.698	-3.6
1,3-Dichlorobenzene	0.779	0.789	-1.3
1,4-Dichlorobenzene	0.788	0.827	-4.9 *
Benzyl Alcohol	0.372	0.4670	
1,2-Dichlorobenzene	0.735	0.761	-3.5
2-Methylphenol	0.547		-7.3
bis(2-Chloroisopropyl)Ether	1.029		-31.6
4-Methylphenol	0.621		
N-Nitroso-di-n-propylamine_		l .	-27.8 #
Hexachloroethane	0.307	1	
Nitrobenzene	0.372		
Isophorone	0.728		
Z NICIOPHENOI	• 0.238	1	
2,4-Dimethylphenol	0.320	4	7.8
Benzoic Acid	0.067	1 .	
bis(2-Chloroethoxy)Methane_	0.507	1	
2,4-Dichlorophenol	0.357	0.315	1 .
1,2,4-Trichlorobenzene	0.392	0.349	11.0
Naphthalene	1.013		1
4-Chloroaniline	0.447	0.469	, ,
	* 0.208		3
4-Chloro-3-methylphenol	* 0.321	0.308	
2-Methylnaphthalene	0.687		
,	# 0.383	•	1
	* 0.432		
2,4,5-Trichlorophenol	0.396		1
2-Chloronaphthalene	1.113	4	1 1
2-Nitroaniline	0.352		1 1
Dimethyl Phthalate	1.206		1
Acenaphthylene	1.666	1.792	
2,6-Dinitrotoluene	0.329		1. 1
3-Nitroaniline	0.329		
	* 0.946		
	# 0.189		
4-Nitrophenol	# 0.071	0.083	-16.9 #
	1		

7A VOLATILE CONTINUING CALIBRATION CHECK

 Lab Name:
 CH2M HILL/LRD
 Contract:
 V28909

 Lab Code:
 Case No.:
 V28909
 SAS No.:
 SDG No.:
 GC-MS

 Instrument ID:
 5100
 Calibration date:
 03/01/91
 Time:
 0911

Lab File ID: 91M2V01683 Init. Calib. Date(s): 02/22/91 02/22/91

Matrix:(soil/water) WATER Level:(low/med) LOW Column:(pack/cap)

Min RRF50 for SPCC(#) = 0.300 (0.250 for Bromoform) Max %D for CCC(#) = 25.0%

COMPOUND	RRF	RRF50	%D
	1.579	1.347	1 1
Bromomethane	1.262	1.293	, ,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,
	1.337	1.315	Ţ
Chloroethane	0.604	0.593	
Methylene Chloride	1.741	2.014	
Acetone	1.326		1 1
Carbon Disulfide	3.504		
1,1-Dichloroethene *		1.029	
	2.780	3.045	
1,2-Dichloroethene (total)	1.587	1.897	i :
Chloroform *	3.063	3.808	
1,2-Dichloroethane	2.296		=35.3
2-Butanone (MEK)	0.093		-21.5
1,1,1-Trichloroethane	0.398	0.419	
Carbon Tetrachloride	0.487		-16.2
Vinyl Acetate	0.846	1	1
Bromodichloromethane	0.682		-32.8
(.)	0.428		-19.6 *
cis-1,3-Dichloropropene	0.533	0.660	
Trichloroethene	0.387	0.496	
Dibromochloromethane	0.593		=53.5
1,1,2-Trichloroethane	0.351	0.527	
Benzene	0.928	1.089	
trans-1,3-Dichloropropene	0.527	7	-33.0
	0.555	0.894त	-61.1
4-Methyl-2-Pentanone	1.225	1.631	-33.1
2-Hexanone	1.086	1.501	-38.2
Tetrachloroethene	0.397	0.517	-30.2
1,1,2,2-Tetrachloroethane #	0.943	1.324	-40.4
Toluene *	i i	0.856	
Chlorobenzene #	0.916	1.172	
	0.424	0.530	-25.0 *
Styrene	0.835	1.103	-32.1
Xylenes (total)	0.489	0.635	-29.9
	======	=======	
Toluene-d8 - SS	0.563	0.523	7.1
1,4-Bromofluorobenzene - SS	0.401	0.393	
1,2-Dichloroethane-d4 - SS	1.026	1.092	
			-

CASE: 28909 SITE: BKGD ANALYSIS: INORG LAB NOTES: -001 diluted Pb in GFAA 1:10 to get reading	
LAB NOTES: -001 diluted PD in CFAA 1:10 to get reading	
Lab reported Pb by ICP analysis for sample 1	
HOLDING TIMES: No analysis time listed. (I used prep date on the spread sheet	
NUMBER OF SAMPLES: Soil 7 Water /	•
CALIBRATION Initial: of	
Continuing: ok	
CONTESTION ICS-OK COEFFICIENT: Not provided ICS-OK	
CONTENT: Not provided ICS-ok COEFFICIENT: Not provided ICS-ok Aqueous - Sb - 8.2 Ba 2.6 Pb 1.0 Zn 4.2 BLANKS Method: Soil Cu 11.1 Ag 1.4	
Field: Cu 2.4 2n 5.0; Flag all adjusted concentrations below 25 in acc	d Gu
FIELD DUPLICATES: -907 + -008 - good match	
SURROGATES:	
MS/MSD: Pre-spike out because conc. too low for He, + Se - no action	
INTERNAL STANDARDS: NA TUNING: NA	
OTHER: LCS - ok	
SUMMARY: The CRDL for lead of 3 and for As of 10 could not be	
met when analysis was by ICP for both (IDL of 13 for As and	
18 for Pb)	

CASE: 28922 SITE: BK&D ANALYSIS: VOC
LAB NOTES: MeCl + Acetone in blank
MS/MSD with 29017-004
HOLDING TIMES:
NUMBER OF SAMPLES: Soil 3 Water 1 ERB - No trip blen
CALIBRATION Initial: MEK RRF of 0.090 - ok for Hazwrep (im.t = 0.05)
Continuing: OK
COEFFICIENT:
BLANKS Method: Not done for water . Soil = MeCl 4 Acetone = 2
Field: " 4 " 25
FIELD DUPLICATES: None indicated
SURROGATES:
MS/MSD: OK
INTERNAL STANDARDS: ok TUNING: ok
OTHER:
SUMMARY: There should have been a method blank for the water sample -
I can't find any done on that date
The soil blank accorded with this case was dated a daty earlier -
that may be a midnight analysis
Nothing besides blank contempation found above CRDL

	5.00
CASE: 28922 SITE: BK6-D	_ ANALYSIS:
LAB NOTES: Blank contam acceptable	
MS/msD with 29017.004	
HOLDING TIMES: W. 3/4 - 2/27 ok Soil	2/28 - 2/27 ok
NUMBER OF SAMPLES: Soil3	Water/
CALIBRATION Initial: OK	
	049
Continuing: RRF for Benzoic Acid: 0.024 HxCCP Diene 36.8 900 = 6	\$ 8D BCIE 36.9
HxCCP Diene 36.8 500 = 6	4.2 Nitropul de 36.6
COEFFICIENT:	plenol 27.5 Benzo (ghi)p. 57.6
BLANKS in - nothing Soil. BEHP 61 EJ a Method: Neither blank-listed was analyzed w Field:	nd diburyi phthelia 43BJ 450 y botch but they were extract w/ south
FIELD DUPLICATES: none	
SURROGATES: PHL-d5 106 out of 10-94 - only 1	so as aclus
MS/MSD: ok	
INTERNAL STANDARDS: ok TUN	ING: 0k
OTHER:	
SUMMARY: No hits associated by CColib.,	no flaga.

CASE: 28922 SITE: BKGD ANALYSIS: /NORG
LAB NOTES: All blanks <crdl; acceptable<="" but="" ke="" ke,="" out="" pb="" se="" sp.="" td=""></crdl;>
RPDs of As + Zn out - LCS varied for furnace but acceptable
ms/msp in 28909
HOLDING TIMES:
NUMBER OF SAMPLES: Soil 3 Water 1
CALIBRATION Initial: ok
Continuing: ok LCS: oK
COEFFICIENT: 1C5 oK Ag = 56 8.2; Ba 2.6; Pb 1.0; Zn 4.2 BLANKS Method: 50.1 = Ca 11.1; Ag 1.4 Field: Ba 3.7; Cu 72.1 Pb 21.9 Zn 50.1
FIELD DUPLICATES: No field dupo
SURROGATES:
MS/MSD: ∘<
INTERNAL STANDARDS: TUNING:
OTHER: As analyzed by ICP - no explanation why in lab notes
OTHER: As analyzed by ICP - no explanation why in lab notes since there were concentrations reported for all samples, it probably doesn't matter
SUMMARY:
·

29034 2 +/ 20224575. 1/00
CASE: 28934 SITE: $2+1$ ANALYSIS: VOQ
LAB NOTES: The 1 ST analysis for med level had invalid IC - The ner-analysis was past holding times (2) Me CI + acetone in blanks
ms/msD with 29017
HOLDING TIMES: 3/13 analysis - 2/28 sampled - 13 days for low
HOLDING TIMES: 3/13 analysis - 2/28 sampled - 13 days for low 1st dilution - 14 days from extraction; 2nd dil - 18 days from extraction NUMBER OF SAMPLES: Soil 3 Water
Initial: Citionathone 35670 RSD of
Continuing: 1,1,2-TCA 30.8 % D chloromethere 27.6% D, 28.6%
Me Cl - 99.9 % D
COEFFICIENT:
BLANKS Method: McCl 4, 10 Acctone 6, 5
Field: none
FIELD DUPLICATES: none
SURROGATES: control charts show % & within range - ok
INTERNAL STANDARDS: * TUNING: de .
INTERNAL STANDARDS: K TUNING: K
OTHER:
SUMMARY: I used the concertration for the re-analysis that lad a good
calibration but was past holding times (24,000 ug/kg) rather than the
invalid IC within acceptable " " (29,000 ug/kg). all other hits recorded
from orizinal analysis.

ASE: 28934 SITE: 2	ANALYSIS: 510C
AB NOTES: blank contam. Ms/MSD	w 29017

OLDING TIMES: 3/4 extract - 2/	28 sampling = 4 days
TUMBER OF SAMPLES: Soil3	Water
CALIBRATION Initial: +k	
Continuing: Benzoic Acid 0.049 RRFS	50 × 26.9 %D
COEFFICIENT:	
BLANKS Method: N-nitrosodiphenylanie (35)	+ BENP (34)
Field: none	
FIELD DUPLICATES: none	
SURROGATES: control charts in	
MS/MSD:	
INTERNAL STANDARDS: %	TUNING: &
OTHER:	
CITMMADV •	
SUMMARY:	
SUMMARY:	
SUMMARY:	

LAB NOTES: Oblank contan, (2) Se post out, 96 preds out, dups out (3) LCS HOLDING TIMES: Hg 3/12 prep - 2/28 samp - 12 days. NUMBER OF SAMPLES: Soil / Water CALIBRATION Initial: ok Continuing: ok Continuing: ok Coefficient: no furnace data supplied BLANKS Method: Ag 1.4 Ca 11.1 Field: none FIELD DUPLICATES: none ; lab dup form not included SURROGATES: MS/MSD: INTERNAL STANDARDS:	CASE: 28934 SITE: 2 ANALYSIS: MORG
HOLDING TIMES: Hg 3/12 prep - 2/28 samp - 12 daya NUMBER OF SAMPLES: Soil / Water CALIBRATION Initial: ok Continuing: ok Continuing: ok Continuing: ok Coefficient: no furnace data supplied BLANKS Method: Az 1.4 Ca 11.1 Field: none FIELD DUPLICATES: none ; lab dup form not included SURROGATES: MS/MSD: INTERNAL STANDARDS: TUNING: OTHER: 1C5 - ok LCS - ok SUMMARY: I couldn't find post-digestion results for furnace metals This lab does not give analysis dates for metals - I used prep dates Forms missing - 5 R + 5 B (pre+post d. spike) 6 (duplicates)	
NUMBER OF SAMPLES: Soil Water CALIBRATION Initial: ok Continuing: ok Continuing: ok Coefficient: no furnace data supplied BLANKS Method: Az 1.4 Cu 11.1 Field: none Jab dup form not included SURROGATES: MS/MSD: TUNING: INTERNAL STANDARDS: TUNING: OTHER: 1C5 - ok LCS - ok LCS - ok SUMMARY: I couldn't find post-digestion results for furnace metals This lab does not give analysis dates for metals - t used prep dates Forms missing - 5A + 5B (pre+post d. spike) 6 (duplicates)	
NUMBER OF SAMPLES: Soil Water CALIBRATION Initial: ok Continuing: ok Continuing: ok Coefficient: no furnace data supplied BLANKS Method: Az 1.4 Cu 11.1 Field: none Jab dup form not included SURROGATES: MS/MSD: TUNING: INTERNAL STANDARDS: TUNING: OTHER: 1C5 - ok LCS - ok LCS - ok SUMMARY: I couldn't find post-digestion results for furnace metals This lab does not give analysis dates for metals - t used prep dates Forms missing - 5A + 5B (pre+post d. spike) 6 (duplicates)	
NUMBER OF SAMPLES: Soil Water CALIBRATION Initial: ok Continuing: ok Continuing: ok Coefficient: no furnace data supplied BLANKS Method: Az 1.4 Cu 11.1 Field: none Jab dup form not included SURROGATES: MS/MSD: TUNING: INTERNAL STANDARDS: TUNING: OTHER: 1C5 - ok LCS - ok LCS - ok SUMMARY: I couldn't find post-digestion results for furnace metals This lab does not give analysis dates for metals - t used prep dates Forms missing - 5A + 5B (pre+post d. spike) 6 (duplicates)	HOLDING TIMES: Hg 3/12 prep - 2/28 samp - 12 days
Continuing: ok Coefficient: no furnace data supplied BLANKS Method: As 1.4 Ca 11.1 Field: none FIELD DUPLICATES: none ; lab dup form not included SURROGATES: MS/MSD: INTERNAL STANDARDS: TUNING: OTHER: 1C5 - ok LCS - ok SUMMARY: I couldn't find post-digestion results for furnace metals This lab does not give analysis dates for metals - I used prep dates Forms missing - 5A + 5B (pre+post d. spike) 6 (duplicates)	•
COEFFICIENT: no furnace data supplied BLANKS Method: As 1.4 Ca 11.1 Field: none FIELD DUPLICATES: none; lab dup form not included SURROGATES: MS/MSD: INTERNAL STANDARDS: TUNING: OTHER: 1C5 - ok LCS - ok SUMMARY: I couldn't find post-digestion results for furnace metals This lab does not give analysis dates for metals - I used prep dates Forms missing - 5A + 5B (pre+post d. spike) 6 (duplicates)	CALIBRATION Initial: ok
BLANKS Method: As 1.4 Ca 11.1 Field: none FIELD DUPLICATES: none; lab dup form not included SURROGATES: MS/MSD: INTERNAL STANDARDS: TUNING: OTHER: 1C5 - ok LCS - ok SUMMARY: I couldn't find post-digestion results for furnace metals This lab does not give analysis dates for metals - I used prep dates Forms missing - 5A + 5B (pre+post d. spike) 6 (duplicates)	Continuing: ok
Method: As 1.9 Ca 11.1 Field: none FIELD DUPLICATES: none; lab dup form not included SURROGATES: MS/MSD: INTERNAL STANDARDS: OTHER: 105 - ork LCS - ork SUMMARY: I couldn't find post-digestion results for furnace metals This lab does not give analysis dates for metals - I used prep dates Forms missing - 5A + 5B (pre+post d. spike) 6 (duplicates)	COEFFICIENT: no furnace data supplied
FIELD DUPLICATES: _none ; lab dup form not included SURROGATES: MS/MSD: INTERNAL STANDARDS:	BLANKS Method: Ag 1.4 Ca 11.1
SURROGATES: MS/MSD: INTERNAL STANDARDS: OTHER: 105 - ok LCS - ok SUMMARY: I couldn't find post-digestion results for furnace metals This lab does not give analysis dates for metals - t used prep dates Forms missing - 5A + 5B (pre + post d. spike) 6 (duplicates)	Field: none
INTERNAL STANDARDS: OTHER: 1C5 - ok LC5 - ok SUMMARY: I couldn't find post-digestion results for furnace metals This lab does not give analysis dates for metals - I used prep date: Forms missing - 5A + 5B (pre+post d. spike) 6 (duplicates)	FIELD DUPLICATES: none; lab dup form not included
INTERNAL STANDARDS: OTHER: 1C5 - ok LCS - ok SUMMARY: I couldn't find post-digestion results for furnace metals This lab does not give analysis dates for metals - I used prep date: Forms missing - 5A + 5B (pre + post d. spike) 6 (duplicates)	SURROGATES:
OTHER: 105 - ok LCS - ok SUMMARY: I couldn't find post-digestion results for furnace metals This lab does not give analysis dates for metals - I used prep date: Forms missing - 5A + 5B (pre + post d. spike) 6 (duplicates)	MS/MSD:
SUMMARY: I couldn't find post-digestion results for furnace metals This lab does not give analysis dates for metals - I used prep dates Forms missing - 5A + 5B (pre + post d. spike) 6 (duplicates)	INTERNAL STANDARDS:TUNING:
SUMMARY: I couldn't find post-digestion results for furnace metals This lab does not give analysis dates for metals - I used prep dates Forms missing - 5A + 5B (pre + post d. spike) 6 (duplicates)	OTHER: 1C5 - ok
This lab does not give analysis dates for metals - I used prep date: Forms missing - 5A + 5B (pre + post d. spike) 6 (duplicates)	
Forms missing - 5A + 5B (pre + post d. spike) 6 (duplicates)	SUMMARY: I couldn't find post-digestion results for furnace metals
Forms missing - 5A + 5B (pre + post d. spike) 6 (duplicates)	This lab does not give analysis dates for metals - I used prep date
6 (duplicates)	·
Ala cartal charts	·
ing control curris	No control charts

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CASE: 28934 SITE: 2 ANALYSIS: BTEX
LAB NOTES: Med level analysis for 8, 9, 11, 12 due to high PAH background
LAB NOTES: Med level analysis for 8, 9, 11, 12 due to high PAH background 5x dilution for -11 +12 due to interference (3) r ² ~ 0.995 for confirmation only
4 = manager To, low for -11
anely. 10w estact HOLDING TIMES: 3/07 - 2/28 = 7 days + 3/11 - 2/28 = 11 days for mediu
NUMBER OF SAMPLES: Soil 10 Water 2 = 1 ERB + 1 TRIP
CALIBRATION Initial: ** for BTEX compounds
Continuing: ok for "
COEFFICIENT: original - ok configuration - < .995 BLANKS Method: no hits
Field: no hits
FIELD DUPLICATES:
SURROGATES: 6990R for - 11 but this sample showed highest concentrations - "J" 5590R for - 10 - not mentioned in lab notes - 1 small hit - UJ for non-
MS/MSD: 6k
INTERNAL STANDARDS: TUNING:
OTHER: control chart - within range
SUMMARY: low surrogate R for 2 samples

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

02B-S-12-14

ŕ	Lab Name: CH2M HILL/LRD Contract:	: <u>S28934 </u> _	
	Lab Code: Case No.: <u>S28934</u> SAS No.:	: SDG N	No.: <u>GC-MS</u>
	Matrix: (soil/water) SOIL	Lab Sample ID:	28934002
4	Sample wt/vol: (g/mL) G	Lab File ID:	91M1BN1784
	Level: (low/med) LOW	Date Received:	03/01/91
	% Moisture: not dec dec	Date Extracted:	03/04/91
3	Extraction: (SepF/Cont/Sonc) SONC	Date Analyzed:	03/23/91
	GPC Cleanup: (Y/N) N pH:	Dilution Factor	: 1

Number TICs found: 11

CONCENTRATION UNITS: (ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q =====
1. 2. 556-67-2 3. 57-10-3 4. 5. 6. 238-84-6 7. 123-79-5 8. 27208-37-3 9. 205-82-3 10. 11.	UNKNOWN CYCLOTETRASILOXANE, OCTAMETH HEXADECANOIC ACID UNKNOWN UNKNOWN 11H-BENZO[A] FLUORENE HEXANEDIOIC ACID, DIOCTYL ES CYCLOPENTA[CD] PYRENE BENZO[J] FLUORANTHENE UNKNOWN UNKNOWN	6.87 9.85 23.07 23.19 24.39 26.31 27.59 28.29 33.39 40.16 40.56	700 190 590 220 270 220 9600 210 510 190	J BJ J J J J J J J J



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EPA SAMPLE NO.

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET TENTATIVELY IDENTIFIED COMPOUNDS

02B-S-18-20

Lab Name: CH2M HILL/LRD	_ Contract:	S28934	
Lab Code: Case No.: S2893	34 SAS No.:	SDG 1	No.: GC-MS
Matrix: (soil/water) SOIL		Lab Sample ID:	28934003
Sample wt/vol: (g/mL)	G	Lab File ID:	91M1BN1785
Level: (low/med) LOW		Date Received:	03/01/91
% Moisture: not dec. 23 dec.		Date Extracted:	03/04/91
Extraction: (SepF/Cont/Sonc)	SONC	Date Analyzed:	03/23/91
GPC Cleanup: (Y/N) N pH:		Dilution Factor:	: 1

CONCENTRATION UNITS: (ug/L or ug/Kg) UG/KG

Number TICs found: 9

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1. 2. 3. 4. 556-67-2 5. 4032-93-3 6. 872-50-4 7. 57-10-3 8. 9. 103-23-1	UNKNOWN UNKNOWN UNKNOWN CYCLOTETRASILOXANE, OCTAMETH HEPTANE, 2,3,6-TRIMETHYL- 2-PYRROLIDINONE, 1-METHYL- HEXADECANOIC ACID UNKNOWN HEXANEDIOIC ACID, BIS(2-ETHY	10.54 10.84 23.07 24.39	250 190 1000 280 190 180 470 520	J J J BJ J J J

000145

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

02B-S-0-2

,	Lab Name: CH2M HILL/LRD Conti	ract: <u>\$28934</u>	_
		No.: SDG No.: GC-MS_	
	Matrix: (soil/water) SOIL	Lab Sample ID: 28934001	
	Sample wt/vol: (g/mL) G	Lab File ID: 91M1BN1783	
	Level: (low/med) LOW	Date Received: 03/01/91	
	% Moisture: not dec. 22 dec	Date Extracted: 03/04/91	
	Extraction: (SepF/Cont/Sonc) SONC	Date Analyzed: 03/23/91	

GPC Cleanup: (Y/N) N pH: ___ Dilution Factor: 1

CONCENTRATION UNITS: (ug/L or ug/Kg) UG/KG

Number TICs found: 8

CAS NUMBER COMPOUND NAME	RT	EST. CONC.	Q =====
1. 147-85-3	TL- 10.52 2 ACID 22.25 23.07	260 200 1200 200 230 210 330 560	BJ J J J J J

CASE: 28934 SITE: 2 ANALYSIS: MORG
LAB NOTES: % R of Se OUT % D of As, In out LCS variance
ms/msD w/ 28909 -001
HOLDING TIMES: Hy Prep 3/12 - Sampling 2/26 = 14 days
NUMBER OF SAMPLES: Soil/ Water
CALIBRATION
Initial: ok
Continuing: ok
COEFFICIENT: can't find
Method: Cu · 11.1 As 1.4 Flag Cu with B: no Ag present
Field: None
FIELD DUPLICATES: None
SURROGATES:
MS/MSD: No problem
INTERNAL STANDARDS: TUNING: *
OTHER:
summary: No analysis times recorded - I used preptime
SUPPLIES TO CONTROL TO CONTROL TO SEE PREP TIME

CASE: 28972 SITE: ANALYSIS:
LAB NOTES: Sample had to be diluted 1:2
co-elution of indeno (1,2,3-cd) pyrene and dibengo
(a,h) anthracene
HOLDING TIMES: 3/06 extract - 3/4 Sampling = 2 days
NUMBER OF SAMPLES: Soil Water /
CALIBRATION Initial: 76RSD for benzo (9hi) perylene - 30.18; confirm = 7.6
Continuing: 900 for benzo (g, h, i) perylene - 40.63
COEFFICIENT: not found
BLANKS Method:mo hits
Field:
FIELD DUPLICATES:monc
surrogates: ot
MS/MSD: not & requested
INTERNAL STANDARDS: TUNING:
OTHER:
SUMMARY:

cian 20972 arms / warvara Affe	
CASE: 28972 SITE: / ANALYSIS: BTEX	
LAB NOTES: orig. analysis out of nence reanalysis at 5x dile	Tio
deleted sample 3/18-3/4 = 14 days HOLDING TIMES: exceeded for confirmation; orig analysis 3/14-3/4=10d	
HOLDING TIMES: exceeded for confirmation; orig analysis 3/14 - 3/4 = 10d	ays
NUMBER OF SAMPLES: Soil Water / PSBW1	
CALIBRATION Initial: οκ	
Continuing: ok for targel compounds (out for styrene, etc.)	
COEFFICIENT:	
BLANKS Method: no detecto	
Field:	
FIELD DUPLICATES:	
SURROGATES: % & within control chart range	
SURROGATES: 70 R within control chart range lot MS/MSD: Note that acceptance criteria were met but no data	nelye
INTERNAL STANDARDS: TUNING:	reft
OTHER:	
SUMMARY:	

CASE: 28973 SITE: QC	ANALYSIS: VOA
LAB NOTES: Travel Blank had to be d	iluted
Me Cl + Acetone found in	method blank
ms/msd w/ 29017.015	
HOLDING TIMES: 3/11 - 3/04 - OK	
NUMBER OF SAMPLES: Soil	
Initial: RRF ok 90RSL ok	IFB; I Trip BI
continuing: Chloromethane - 31.6	7- not found in sample-
	no action t
COEFFICIENT: NA	
Method: MeCland Acetona 12 Field: MeCl Acetona FIELD DUPLICATES: none	
surrogates: ok	
SURROGATES &	
WC (WCD.	
MS/MSD:	runing: 🏒
	runing: 🥕
INTERNAL STANDARDS: OK	runing: 🍂
INTERNAL STANDARDS: OK	runing: 🎻
INTERNAL STANDARDS: <u>ok</u>	runing: 🎤
INTERNAL STANDARDS: <u>ok</u>	runing: <u>ok</u>
INTERNAL STANDARDS: <u>ok</u>	runing: 🎤

CASE: 28973 SITE: QC ANALYSIS: SVOC
LAB NOTES: some blank contam. but within critoria
ms/msD with 29017-015
HOLDING TIMES: 3/04 est - 3/04 - 0K
NUMBER OF SAMPLES: Soil Water _/ FB
CALIBRATION
Continuing: Benzoic Acid47.8 %); Benzo [ghi] parybu -43.8%
Hexachloro futadienes cyclopentadienes 29%D
- COEFFICIENT:
BLANKS Method: both compounds found in FB were found in method blank
Field: the I semple is a field blank
FIELD DUPLICATES: none indicated
SURROGATES: ok
MS/MSD:
INTERNAL STANDARDS: TUNING: *
OTHER:
SUMMARY:

CASE: 28973 SITE: 9C	ANALYSIS: /NORG
LAB NOTES: DUP + SPIKE Wy 29017.	tor all but Hg
2890	9.010 for Hg
•	
HOLDING TIMES: No date of analysis 50	rea - ok using prepdele
NUMBER OF SAMPLES: Soil	Water/
CALIBRATION Initial: ok	
Continuing: ok	
COEFFICIENT: not found	
BLANKS Method: 5.98 Zo	
Field: This one sample is a field A	lank - has a copper hit of 51.
FIELD DUPLICATES: None	
SURROGATES:	
MS/MSD:	
INTERNAL STANDARDS:	TUNING: LCS: OK
OTHER:	
SUMMARY:	
•	
,	

CASE: 28998 SITE: 1 ANALYSIS: BTEX
LAB NOTES: Holding times exceeded for 2 samples and the ms/msD run009 and -011 were diluted because of matrix interference
additional information attached
HOLDING TIMES: Analysis 3/15 · 20 - Sampling date 3/5 = [15] days
HOLDING TIMES: Analysis 3/15-20 - Sampling date 3/5= 15 days NUMBER OF SAMPLES: Soil 16 for -002 and -010 and MS/MSI Water CALIBRATION
CALIBRATION Initial: ok
Continuing: ok
COEFFICIENT: - half are low - all associated concentrations should
BLANKS DE LA II MERCO
Method: nothing found
Field: no detecto
FIELD DUPLICATES: possible -002 + -003 / Thit is one -ND in other
surrogates: 10w 35%R +52%R high 188%R - 206%R
MS/MSD:
INTERNAL STANDARDS: TUNING:
OTHER: They re-ran he high suspogated with Similar results.
didn't re run the low surrogates because There would still
SUMMARY: be an interjerence problem.
Flag non detecta UJ because of low correlation coefficients
and low surrogate recoveries.

CASE: 29017 SITE: 2 ANALYSIS: VOA	
LAB NOTES: -006, -011, -016 were diluted (16 is a trip blank)	
MeCl + Acetone + TCE in method blank	
	L.
ms/msD: TCE out, diluted sample run for spike but not for duplica	ИC
HOLDING TIMES: Analysis 3/20 - Sampling 3/6 - OK	
NUMBER OF SAMPLES: Soil 14 Water 2 -nit. for soils, given are rel for medium level -none given for low level so All CALIBRATIONS Initial: Chloromethane (med soil) 41.8 % RSD MeC1 30.7	il
Acetone 31.6% MEK 0.041 RRF	
Continuing: med soil Chloromethane 33.5% D. MEK 0.038 RRi	=57
CS2 - 30.4 %D	
COEFFICIENT: Low soil - see attached	
BLANKS Method: see attacked	
Field: ERB: Acetone 36; MeCl 4, TCE 1; TRAVEL: CS_1800; M	eCl
FIELD DUPLICATES: -013 + -014 9 -003 + -004 -9000 to thirty 9000	
SURROGATES: OK	1
MS/MSD: ok - 1 out but ok	
INTERNAL STANDARDS: OK TUNING: OK	
OTHER: There were a number of TICs in both medium + low soil	
samples - examples are attached	
SUMMARY:	
	•

CASE: 29017 SITE: 2 ANALYSIS: SVOC
LAB NOTES: - 1 werd; luted blank contan was acceptable
ms/msd, within limits
HOLDING TIMES:
NUMBER OF SAMPLES: Soil /4 Water /
CALIBRATION Initial: OF
Continuing: Berroic Acid -99.9, -91.8, 82.5 Berro 9hi 32.8 Hexcelloro cyclopertadiene, 53.7, 39.6 COEFFICIENT: 3-Nitroaniline 40.5 4-Nitroaniline 47.1 [no Hits, no flag
Hexaclorocyclopertadione, 53-7, 39.6 2.4-diritro phenol 33.7
COEFFICIENT: 3-Nitroaniline 40.5 4-Nitroaniline 47.1 [no Hits, no fla
BLANKS Method: BEHP. 66 + 73 in soil blank - B removed from all over 800
Field: BEHP. 5
FIELD DUPLICATES: good matches
surrogates: Phenol-d5 - high 70R in 4 of 6 samples - I did not flag
MS/MSD: because it was all for a single surrogate, but you may wish to. (Both 90Rs out for 2 compounds - LCS ok - I, took no action
INTERNAL STANDARDS: ok TUNING: OK.
OTHER:
SUMMARY: Lots of TIC, - some examples attached
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CHM HILL TELEPHONE CONVERSATION RECORD

CALL TO		•	······································	PHONE NO.	
CALL FRO				TIME	
MESSAG	E TAKEN BY			PROJECT NO.	
SUBJECT					
Lab	Sample Its	N-nitrosodio.	d: . a . bytul . a	ВбНР	·
	-15, MIE, DIS	3			
w z			~	4	
<u>s1</u>	-1 - -9	-		66	
<i>5</i> 2	-1014, +14	44 We/wad		73	•
				40	
	•				
				-	

1F SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

Lab Name: CH2M HILL/LRD Contract: S29017

Lab Code: _____ Case No.: <u>S29017</u> SAS No.: _____ SDG No.: <u>GC-MS</u>

Matrix: (soil/water) SOIL Lab Sample ID: 29017009

Sample wt/vol: ____ (g/mL) <u>G</u> Lab File ID: <u>91M1BN1859</u>

Level: (low/med) LOW Date Received: 03/07/91

% Moisture: not dec. 21 dec. Date Extracted: 03/11/91

Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 03/29/91

GPC Cleanup: (Y/N) N pH: ____ Dilution Factor: 1

CONCENTRATION UNITS: (ug/L or ug/Kg) UG/KG

Number TICs found: 20

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1.	UNKNOWN	6.90	59000	J
2. 095-47-6	BENZENE, 1,2-DIMETHYL-	7.33	3400	J
3. 13475-81-5	HEXANE, 2,2,3,3-TETRAMETHYL-		6000	J
4. 62016-34-6	OCTANE, 2,3,7-TRIMETHYL-	8.59	7600	J
5. 52896-87-4	HEPTANE, 4-(1-METHYLETHYL)-	8.74	3700	J
6.	UNKNOWN	9.00	3700	J
7. 620-14-4	BENZENE, 1-ETHYL-3-METHYL-	9.17	12000	J
8.	UNKNOWN	9.29	5000	J
9. 124-18-5	DECANE	9.84	20000	J ·
10. 17302-28-2	NONANE, 2,6-DIMETHYL-	10.25	5300	J
11. 108-67-8	BENZENE, 1,3,5-TRIMETHYL-	10.34	4200	J
12.	UNKNOWN	10.47	6200	J
13. 1074-55-1	BENZENE, 1-METHYL-4-PROPYL-	10.82	4300	J
14.	UNKNOWN	10.95	8800	J
15. 62016-37-9	OCTANE, 2,4,6-TRIMETHYL-	11.57	10000	J
16. 767-99-7	BENZENE, (1-METHYL-1-PROPENY	12.55	2900	J
17. 30571-71-2	DECANE, 3-BROMO-	12.74	2000	J
18. 2801-84-5	DECANE, 2,4-DIMETHYL-	14.74	4300	J
19. 31295-56-4	DODECANE, 2,6,11-TRIMETHYL-	16.15	2900	J
20. 74645-98-0	DODECANE, 2,7,10-TRIMETHYL-	17.52	1600	J

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

·Lab	Name:	CH2M HILL/LRD	Contract: S29017	04BS-14-16D
Lab	Code:	Case No.: <u>S29017</u>	SAS No.: SDG	No.: GC-MS

Lab Sample ID: 29017004 Matrix: (soil/water) SOIL

Sample wt/vol: ____ (g/mL) G Lab File ID: 91M1BN1854

Level: (low/med) LOW Date Received: <u>03/07/91</u>

% Moisture: not dec. ___ dec. ___ Date Extracted: 03/11/91

Extraction: (SepF/Cont/Sonc) SONC Date Analyzed: 03/29/91

GPC Cleanup: (Y/N) N pH: ___ Dilution Factor: 1____

CONCENTRATION UNITS: Number TICs found: 20 (ug/L or ug/Kg) <u>UG/KG</u>

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1. 075-91-2	HYDROPEROXIDE, 1,1-DIMETHYLE	6.88	56000	BJ
2. 095-47-6	BENZENE, 1,2-DIMETHYL-	7.30	1100	J
3. 13475-81-5	HEXANE, 2,2,3,3-TETRAMETHYL-	7.90	1200	J
4. 5911-04-6	NONANE, 3-METHYL-	8.57	1900	J
5. 620-14-4	BENZENE, 1-ETHYL-3-METHYL-	9.15	3100	J
6. 526-73-8	BENZENE, 1,2,3-TRIMETHYL-	9.27	1400	J
7. 124-18-5	DECANE	9.80	6100	J
8. 17302-28-2	NONANE, 2,6-DIMETHYL-	10.22	2100	J
9.	UNKNOWN	10.32	1700	J
10.	UNKNOWN	10.44	2700	J
11. 1074-55-1	BENZENE, 1-METHYL-4-PROPYL-	10.80	2400	J
12.	UNKNOWN	10.94	3200	J
13. 13151-34-3	DECANE, 3-METHYL-	11.07	1300	J
14. 17301-32-5	UNDECANE, 4,7-DIMETHYL-	11.57	5200	J :
15.	UNKNOWN	12.55	1100	J
16. 62016-34-6	OCTANE, 2,3,7-TRIMETHYL-	14.30	860	J
17. 17301-30-3	UNDECANE, 3,8-DIMETHYL-	14.72	* 1800	J
18. 7289-40-9	ETHER, HEPTYL HEXYL	16.14	900	J
19. 057-10-3	HEXADECANOIC ACID	22.92	1200	BJ
20. 103-23-1	HEXANEDIOIC ACID, BIS(2-ETHY	27.42	990	J



1F SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

04BS-14-16

Lab Name: CH2M HILL/LRD	_ Contract: <u>S29017</u>	0486 14 10
Lab Code: Case No.: S290	17 SAS No.: SDG	No.: GC-MS
Matrix: (soil/water) <u>SOIL</u>	Lab Sample ID:	29017003
Sample wt/vol: (g/mL)	G Lab File ID:	91M1BN1853
Level: (low/med) <u>LOW</u>	Date Received:	03/07/91
% Moisture: not dec. 23 dec.	Date Extracted:	03/11/91
Extraction: (SepF/Cont/Sonc)	SONC Date Analyzed:	03/29/91

GPC Cleanup: (Y/N) N pH: ___ Dilution Factor: 1

CONCENTRATION UNITS: (ug/L or ug/Kg) UG/KG

Number TICs found: 20

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1. 075-91-2	HYDROPEROXIDE, 1,1-DIMETHYLE	6.87	60000	вл
2. 13475-81-5	HEXANE, 2,2,3,3-TETRAMETHYL-		1200	J
3. 5911-04-6	NONANE, 3-METHYL-	8.57		J
4.	UNKNOWN	9.14	2800	J
5.	UNKNOWN	9.27	1200	J
6. 124-18-5	DECANE	9.82	5700	J
7. 17302-28-2	NONANE, 2,6-DIMETHYL-	10.22	1800	J
8. 095-63-6	BENZENE, 1,2,4-TRIMETHYL-	10.30	1300	J
9. 7058-01-7	CYCLOHEXANE, (1-METHYLPROPYL		1200	J
10. 17302-32-8	NONANE, 3,7-DIMETHYL-	10.52	1100	J
11. 6975-98-0	DECANE, 2-METHYL-	10.95	5000	J
12. 13151-34-3	DECANE, 3-METHYL-	11.07	1200	J
13. 933-98-2	BENZENE, 1-ETHYL-2,3-DIMETHY	11.25	1000	J
14. 62016-37-9	OCTANE, 2,4,6-TRIMETHYL-	11.57	5100	J
15. 4292-92-6	CYCLOHEXANE, PENTYL-	12.22	700	J
16.	UNKNOWN	12.60	1200	J
17. 62016-34-6	OCTANE, 2,3,7-TRIMETHYL-	14.30	" 860	J
18. 62016-37-9	OCTANE, 2,4,6-TRIMETHYL-	14.70	1400	J
19. 62016-37-9	OCTANE, 2,4,6-TRIMETHYL-	16.14	880	J
20. 057-10-3	HEXADECANOIC ACID	22.92	1200	BJ

000308



1-F

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET TENTATIVELY IDENTIFIED COMPOUNDS

EPA SAMPLE NO.

04BS-6-8

Lab Name: CH2M HILL/LRD Co	ontract: <u>\$29017</u>
Lab Code: Case No.: <u>S29017</u>	SAS No.: SDG No.: GC-MS
Matrix: (soil/water) SOIL	Lab Sample ID: 29017002
Sample wt/vol: (g/mL) G	Lab File ID: 91M1BN1852
Level: (low/med) LOW	Date Received: 03/07/91
% Moisture: not dec dec	Date Extracted: 03/11/91
Extraction: (SepF/Cont/Sonc) SONC	Date Analyzed: 03/29/91

GPC Cleanup: (Y/N) N pH: ____

Dilution Factor: 1

Number TICs found: 20

CONCENTRATION UNITS: (ug/L or ug/Kg) <u>UG/KG</u>

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
1. 141-79-7	3-PENTEN-2-ONE, 4-METHYL-	5.72	480	вл
2. 108-21-4	ACETIC ACID, 1-METHYLETHYL E	6.22	370	BJ
3. 075-91-2	HYDROPEROXIDE, 1,1-DIMETHYLE		53000	BJ
4.	UNKNOWN	8.25	550	J
5. 871-83-0	NONANE, 2-METHYL-	9.15	350	J
6. 124-18-5	DECANE	9.82	1000	J .
7. 13475-78-0	HEPTANE, 5-ETHYL-2-METHYL-	10.22	360	J
8. 6975-98-0	DECANE, 2-METHYL-	10.95	930	J
9. 871-83-0	NONANE, 2-METHYL-	11.57	1200	J
10.	UNKNOWN	11.94	290	J
11. 62183-55-5	OCTANE, 3-ETHYL-2,7-DIMETHYL	12.62	270	J
12. 62016-34-6	OCTANE, 2,3,7-TRIMETHYL-	14.32	340	J
13. 17301-30-3	UNDECANE, 3,8-DIMETHYL-	14.72	580	J
14. 7289-40-9	ETHER, HEPTYL HEXYL	16.15	350	J
15. 62238-13-5	DECANE, 2,3,7-TRIMETHYL-	17.00	200	J
16. 62016-37-9	OCTANE, 2,4,6-TRIMETHYL-	17.50	200	J
17. 057-10-3		22.92	* 550	BJ
18.	UNKNOWN	24.95	230	J
19. 4337-65-9	HEXANEDIOIC ACID, MONO(2-ETH	27.39	1000	J
20. 21078-65-9	1-DECANOL, 2-ETHYL-	28.36	410	J

000300

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET TENTATIVELY IDENTIFIED COMPOUNDS

04BS-0-2

Lab Name: CH2M H	IILL/LRD	Contract: S29017	
Lab Code:	Case No.: <u>S2901</u>	7 SAS No.:	SDG No.: GC-MS
Matrix: (soil/wa	ter) <u>SOIL</u>	Lab Sample	ID: 29017001
Sample wt/vol:	(g/mL)	G Lab File II	91M1BN1851_
Level: (low/m	ned) <u>LOW</u>	Date Receiv	red: 03/07/91
% Moisture: not	dec. 22 dec.	Date Extrac	eted: <u>03/11/91</u>
Extraction: (S	SepF/Cont/Sonc)	SONC Date Analyz	ed: <u>03/29/91</u>

GPC Cleanup: (Y/N) N pH: ___ Dilution Factor: 1

Number TICs found: 20

CONCENTRATION UNITS: (ug/L or ug/Kg) UG/KG

CAS NUMBER	COMPOUND NAME	RT	EST. CONC.	Q
CAS NUMBER ===================================	COMPOUND NAME ===================================	6.92 7.97 8.64 9.05 9.24 9.87 9.94 10.34 10.54	EST. CONC.	Q ===== BJ JJJJJJJJJJJJJJJJJJJJJJJJJJJJJ
19. 17301-30-3 20.	UNDECANE, 3,8-DIMETHYL- UNKNOWN	16.20 32.46	13000 23000	J J



CASE: 29017 SITE: 2 ANALYSIS: /NORG
LAB NOTES: Pre-disestion spike met for Sb, Cr, Se but acceptable post
900 Cr dup. out Se used MSA
HOLDING TIMES: Dale Prepped 3/13 - 3/6 : OK
NUMBER OF SAMPLES: Soil
CALIBRATION Initial: well = ok 5 - ok
Continuing: water or 5 - 0K
COEFFICIENT:
Method: Soil Az-0.205 Se 0.162 Water 5.9 Za
Field: Zn 2.68 (15=13) All cone. found in camples were high no flags
FIELD DUPLICATES:
SURROGATES:
MS/MSD:
INTERNAL STANDARDS:TUNING:
OTHER:
SUMMARY: Pre- + Post- dig spike both low for Se - I for the hits, but
I also gave the one non-hit a UT which you may not went, as
we are not flagging Organic non-detects.

010D 2003 2 0TMD 000
CASE: 29032 SITE: QC ANALYSIS: VOA
LAB NOTES: -003 diluted - CS2 problem -> from inadequate water purification sys
Black contain: Aceton, MeCl + TCE
3) Surrogates - 2 out but acceptable (4) ms/msD with 29017.015
HOLDING TIMES:
NUMBER OF SAMPLES: Soil Water $3 = 2 + 6 + 1 \text{ Trip } B$
CALIBRATION Initial: 9K
Continuing: Acetone 34.3% D no other hit effected
"not present n associated sample - no action
COEFFICIENT: NA (2) MeCI & TCE I BLANKS Method: 1 MeCI8; Acetone 10 (3) MeCI 2, Acetone 8 (b) MeCI 8 Ace 7
Field: both are blanks
FIELD DUPLICATES:
SURROGATES: -003 (Travel Blank) high - all positive results = I
MS/MSD: ok
INTERNAL STANDARDS: ok TUNING: ok
OTHER:
SUMMARY:

CASE: 29032 SITE: QC ANALYSIS: SVOC LAB NOTES: (1) N-mitigardiplenglanine in black (2) -002 low purk. 7R (3) MS/MSD with -2907 HOLDING TIMES: 3/12 (extract.) - 3/7 sampled = 0K NUMBER OF SAMPLES: Soil - Water 2 - both field black CALIBRATION Initial: 0K Continuing: Beagain adid 0.029 KF50, 70.1% Nitroaniline 35.8% D - no hits, no flags COEFFICIENT: NA BLANKS Method: n-Nitrosodiplenglanine 2 - not in samples Field: ooth contain 86HP FIELD DUPLICATES: none SURROGATES: Phead-05 - 476 K - all pointries in sample 2 flagged J DUX + HAZWERP samp non-diffects should be rejected, but limits MS/MSD1 on 10-94 - 2 didn't reject because you didn't havour report And become we don't list all all compounds INTERNAL STANDARDS: 0K TUNING: 0K: OTHER: alkyl groups in TICs 446 - all positives flagged J SUMMARY:	
(2) -002 low purp. TR (3) ms/ms) inth 29017 HOLDING TIMES: 3/12 (extrect.) - 3/7 sempled = 0K NUMBER OF SAMPLES: Soil - Water 2 - both field black CALIBRATION Initial: 0K Continuing: Bengoic adid 0.029 18750, 70.170) Nitroaviline 35.870 - no hits, no flags COEFFICIENT: NA BLANKS Method: n-Nitrosodiphenglamine 2 - not in semples Field: ooth contain BEHP FIELD DUPLICATES: none SURROGATES: Phenol - 45 - 476 k - sel positives in sample 2 flagged J DOWN + HAZWERP samp non-diffects should be rejected but limits METHODE and 10-94-3 disht reject because you disht he your report and because we don't list sele the compounds. INTERNAL STANDARDS: OK TUNING: OK OTHER: alkyl groups in TIC: 476 k - sel positives flagged J	CASE: 29032 SITE: QC ANALYSIS: SVOC
(2) -002 low purp. TR (3) ms/ms) inth 29017 HOLDING TIMES: 3/12 (extrect.) - 3/7 sempled = 0K NUMBER OF SAMPLES: Soil - Water 2 - both field black CALIBRATION Initial: 0K Continuing: Bengoic adid 0.029 18750, 70.170) Nitroaviline 35.870 - no hits, no flags COEFFICIENT: NA BLANKS Method: n-Nitrosodiphenglamine 2 - not in semples Field: ooth contain BEHP FIELD DUPLICATES: none SURROGATES: Phenol - 45 - 476 k - sel positives in sample 2 flagged J DOWN + HAZWERP samp non-diffects should be rejected but limits METHODE and 10-94-3 disht reject because you disht he your report and because we don't list sele the compounds. INTERNAL STANDARDS: OK TUNING: OK OTHER: alkyl groups in TIC: 476 k - sel positives flagged J	LAB NOTES: (N- mitrosodipheny Lanine in black
(3) MS/MSD with 29927 HOLDING TIMES: 3/12 (extrect.) - 3/7 sampled = 0K NUMBER OF SAMPLES: Soil - Water 2 - both field blacks CALIBRATION Initial: 0K Continuing: Bengaic add 0.029 RF50, 70.1% Nitroaviline 35.8% COEFFICIENT: NA BLANKS Method: n-Nitrosodiphenglamine 2 - not in samples Field: ooth contain BEHP FIELD DUPLICATES: none SURROGATES: Phenol-d5 - 4% K- all positives in sample 2 flagged J DOWN + HAZWERP Surpa non-altecto should be rejected, but limits ME/MEB+ are 10-94-2 dioalt reject because you didn't he your report and because we don't list all the compounds. INTERNAL STANDARDS: OK TUNING: 0K OTHER: allyl groups in TIC: 4% R- all positives flagged J	· ·
HOLDING TIMES: 3/12 (extract.) - 3/7 sampled = 0K NUMBER OF SAMPLES: SOII - Water 2 - both field black CALIBRATION Initial: 0K Continuing: Beagain add 0.029 RF50, 70.1%) Nitroanline 35.8% D - no hits, no flags COEFFICIENT: NA BLANKS Method: n-Nitrosodiphenglarine 2 - not in samples Field: ooth contain BEHP FIELD DUPLICATES: none SURROGATES: Phonof - d5 - 4% R - all positives in sample 2 flagged J WY + HAZWEAP Soura mon-dittets should be rejected, but limits MS/MCB1 are 10-94-2 diolet reject because you didn't in your report and because we don't list all the compounds. INTERNAL STANDARDS: 0K TUNING: 0K OTHER: allyl groups in TICs 4% R - all positives flagged J	(3) ms/msD with -015
Continuing: Bengoic alid 0.029 RF50, 70.1% Nitroaniline 35.8%) — no hits, no flags COEFFICIENT: NA BLANKS Method: n-Nitrosodiphenglanine 2 — not in samples Field: ooth contain BEHP FIELD DUPLICATES: none SURROGATES: Phenol-d5 - 4% R-all positives in sample 2 flagged J W + HAZWEAP Soupe mon-diffects should be rejected, but limits MS/MSB: are 10-94-2 docht reject because you didn't in your report and because we don't list all the compounds. INTERNAL STANDARDS: OK TUNING: OK OTHER: alkyl groups in TICs 4% R - all positives flagged J	
Continuing: Bengoic alid 0.029 REFSO, 70.1% Nitroaniline 35.8% D — no hits, no flags COEFFICIENT: NA BLANKS Method: n-Nitrosodiphenglanine 2 — not in samples Field: ooth contain 86HP FIELD DUPLICATES: none SURROGATES: Phenol-d5 - 4% R-all positives in sample 2 flagged J W + HAZWEAP Saya mon-ditecto should be rejected, but limits ME/MEB: are 10-94-2 doubt reject because you didn't in your report and because we don't list all the compounds. INTERNAL STANDARDS: OK TUNING: OK OTHER: alkyl groups in TICs 4% R - all positives flagged J	NUMBER OF SAMPLES: Soil - Water 2 - both field blesk
Nitroaniline 35.8 % D — no hits, no flags COEFFICIENT: NA BLANKS Method: n-Nitrosodiphenglanine 2 — not in samples Field: ooth contain BEHP FIELD DUPLICATES: none SURROGATES: Phenol-d5 — 4% R - all positives in sample 2 flagged J ON + HAZWEAP says mon-diffects should be rejected, but limits MS/MSB: are 10-94 — 2 diolett reject because you didn't in your report and because we don't list all the compounds. INTERNAL STANDARDS: OK TUNING: oK OTHER: alkyl groups in TICs 4% R - all positives flagged J	CALIBRATION
Nitroaniline 35.8 % D — no hits, no flags COEFFICIENT: NA BLANKS Method: n-Nitrosodiphenglanine 2 — not in samples Field: ooth contain BEHP FIELD DUPLICATES: none SURROGATES: Phenol-d5 - 4% R - all positives in sample 2 flagged J ON + HAZWEAP says mon-diffects should be rejected, but limits MS/MSB: are 10-94 - 2 diolate reject because you didn't in your report INTERNAL STANDARDS: OK TUNING: OK OTHER: alkyl groups in TICs 4% R - all positives flagged J	Continuing: Bengair alid 0.029 REF50, 70.1%)
BLANKS Method: n-Nitrosodiphenglanine 2 - not in samples Field: ooth contain BEHP FIELD DUPLICATES: none SURROGATES: Phenol- 45 - 4% R - all positives in sample 2 flagged J ON + HAZWEAP says non-diffects should be rejected, but limits MS/MSB: are 10-94 - 2 diolit reject because you didn't in your report and because we don't list all the compounds. INTERNAL STANDARDS: OK TUNING: oK OTHER: alkyl groups in TIC: 4% R - all positives flagged J	
Method: n-Nitrosodiphenylamine 2 - not in samples Field: poth contain BEHP FIELD DUPLICATES: none SURROGATES: Phenol-d5 - 4% K- all positives in sample 2 flagged J OW + HAZWEAP says non-ditlects should be rejected, but limits MS/MSB: are 10-94 - 2 dight reject because you dight in your report and because we don't list all the compounds: INTERNAL STANDARDS: OK TUNING: OK OTHER: alkyl groups in TICs 4% K - all positives flagged J	COEFFICIENT: NA
SURROGATES: Phenol-d5 - 4% R-all positives in sample 2 flagged J W + HAZWEAP says non-ditects should be rejected, but limits MS/MSB: are 10-94 - 2 dishit reject because you didn't in your report and because we don't list all the compounds. INTERNAL STANDARDS: OK TUNING: oK OTHER: alkyl groups in TICs 4% R - all positives flagged J	BLANKS Method: n-Nitrosodiphenglamine 2 - not in samples
SURROGATES: Phenol-d5 - 4% R-all positives in sample 2 flagged J SW + HAZWEAP says non-attects should be rejected, but limits MS/MSB: are 10-94 - I didn't reject because you didn't in your report and because we don't list all the compounds: INTERNAL STANDARDS: OK TUNING: OK OTHER: alkyl groups in TICs 4% R - all positives flagged J	Field: ooth contain BEHP
TUNING: alkyl groups in TICs 496R - all positives flagged J	FIELD DUPLICATES: none
MS/MSB: are 10-94-S didn't reject because you didn't in your report and because we don't list all the compounds. INTERNAL STANDARDS: OK TUNING: OK OTHER: alkyl groups in TICs 4% R - all positives flagged J	SURROGATES: Phenol-d5 - 4% R-all positives in sample 2 flagged J
OTHER: alkyl groups in TICs 496R - all positives flagged J	OW + HAZWRAP Says non-ditects should be rejected, but limits MS/MSB: are 10-94-2 didn't reject because you didn't in your re
	INTERNAL STANDARDS: OK TUNING: OK
SUMMARY:	OTHER: alkyl groups in TICs 490R - all positives flagged J
SUMMARY:	
	SUMMARY:
"	

	S:	· · · · · · · · · · · · · · · · · · ·			· ·
HOLDING 1	PIMES: 3/19-12	3 — 3/7 sa	mpling =	9K	
NUMBER OF	SAMPLES:	Soil	w	ater <u>2</u>	-
CALIBRATI Initia	on 1: <u>0K</u>	·			
Contir	uing: oK				
ره براد COEFFICII	Cion ENT: Not four	nd			
BLANKS Method	1: Ag -1.2;	Zn 5.9	·		
	Q 11 lan	are field bla	nks both	contain P	6 + Zn
Field	Doth samples		,		
	PLICATES: n				
FIELD DUE	PLICATES:n				
FIELD DUE	PLICATES:n	one			
FIELD DUE SURROGATE MS/MSD:	PLICATES:n	one			
FIELD DUE SURROGATE MS/MSD: INTERNAL	PLICATES:n	ione			
FIELD DUE SURROGATE MS/MSD: INTERNAL	PLICATES:n	ione			
FIELD DUE SURROGATE MS/MSD: INTERNAL OTHER:	PLICATES:n	ione			
FIELD DUE SURROGATE MS/MSD:	PLICATES:n	ione			
FIELD DUE SURROGATE MS/MSD: INTERNAL OTHER:	PLICATES:n	ione			

CASE: <u>29032</u> SITE: <u>4</u>	ANALYSIS: BTEX
LAB NOTES: MJ/MSD W/ 28998-2	<u>.</u>
HOLDING TIMES: 3/21 analysis date -	3/09 sampling date : 14 days
HOLDING TIMES: 3/21 enelyis date -	Water 6 * (2 missing?
CALIBRATION Initial: ok	
Continuing:	
COEFFICIENT: NA	
BLANKS Method: nothing found	
Field: not enalyzed	
FIELD DUPLICATES: not requested	
FIELD DUPLICATES: not requested SURROGATES: ok high for semples 133	5-143; ok for 602/8020-96-11
MS/MSD:	
INTERNAL STANDARDS:	TUNING: Charts . * ok
as samples (note: See ref #) but	list ERB-06-37 and P7BW
as samples (note : See ref # ") but	no mention of these by lab
SUMMARY: no delecto	

CASE: 29032 SITE: 4 ANALYSIS: PNA
LAB NOTES: control charto incomplete
coelution between indeno (1, 2, 3-cd) pyrene and divergo (a, h)entheace
HOLDING TIMES: 3/13 extraction date - 3/7 = 6 days
NUMBER OF SAMPLES: Soil Water
CALIBRATION Initial: high RSD in the metyl, nephtlalenes (41.4 + 38.8) in confirm. Colo
Continuing: 37.3% D for indeno/disenso (coeluting see abore leb
COEFFICIENT:
BLANKS Method: nothing found
Field: Samples 9 +10 Not analyzed
FIELD DUPLICATES: panples 9 + 10
SURROGATES: low for deca 48-59 70 R; low to ok for texp 62-130
MS/MSD: range is 70 - 130; duplicate parge 64-48 61-94
INTERNAL STANDARDS:TUNING:
OTHER:
SUMMARY: nothing found in any samples, so I'm worried about
The low surrogate resorries (can't find acceptable limits
in the Fed Register for 610



POLYNUCLEAR AROMATIC HYDROCARBONS SURROGATE RECOVERY

Primary Analysis Instrument ID: GC 3700

	Date	Laboratory Reference No.	EPA-610 Surrogate Recovery Decafluorobiphenyl	EPA-610 Surrogate Recovery Terphenyl-d14
01	04-04-91	Method Blank	83	-60
02	03-27-91	29032 - 4	49	111
i 03	03-27-91	29032 - 5	48	119
02	03-27-91	29032 - 6	į 54 į	130
03	03-27-91	29032 - 7	57	117
104	03-27-91	29032 - 8	48	121
05	03-27-91	29032 - 9	53	129
06	03-27-91	29032 - 10	58	124
07	03-27-91	29032 - M10	59	62
08	03-27-91	29032 - 010	56	66

EPA-610 Surrogate standards reported as percent recovery.

NA = Not Analyzed.

Comments:

Approved By:

kdh.I

CASE: 29041 SITE: 1 ANALYSIS: BTEX
LAB NOTES: The soil sample at 5x dilution - interference with IS
M5/MSD is acceptable - not included in package
HOLDING TIMES: 3/14 - 3/7 = 7 days
NUMBER OF SAMPLES: Soil / Water / SERB
CALIBRATION Initial: ok
Continuing: ok
COEFFICIENT: not included
BLANKS Method: Clan 112% sum. R
Field: nothing found
FIELD DUPLICATES:
SURROGATES: 6k
MS/MSD: Not available
INTERNAL STANDARDS: TUNING: Control chart ok
OTHER:
SUMMARY:

	#
case: 2904/ site: 1 analysis: PNA	
LAB NOTES: discreparcy between 1st run (130 mg/c) and confirmation	
run (9 ugh) for 1-methyl naphthalere - maticy interference	
HOLDING TIMES:	
NUMBER OF SAMPLES: Soil Water _/	
CALIBRATION Initial: 30.18 % RSD for benzo (9hi) peryline	
I cont tell if IC n CC is for the water on the soil sample	
Continuing:	
COEFFICIENT:	
BLANKS Method: mothers found	
Field: ERB - no lito but surrogete recoveries just above lineto (4	7
FIELD DUPLICATES:	
SURROGATES: 70R for method blank below 43-116 limits = 41	
MS/MSD:	
INTERNAL STANDARDS:	
OTHER: see notes - I gave the 130 a J	
SUMMARY:	

case: 29046 site: 4 analysis: BTEX
LAB NOTES: IS 1000 Surrogates - high ms/msD w/ this package
HOLDING TIMES: 3/21 enalysis - 3/8 sampled = 13 days
NUMBER OF SAMPLES: Soil 6 Water 2 = 1 ERB + 1 Trip
CALIBRATION Initial: dale giren = 2/21
3/15 for standards = ok
Continuing: ok
COEFFICIENT: - date given: 4/1/91 - all ok for that date
BLANKS Method: no lits
Field:mo_iito
FIELD DUPLICATES: no hats
SURROGATES: Lowest Surrogate recovery = 1127 (see attacked) up to 285 % R
MS/MSD: Benzene 182% R dup 265% R Toluene 123 167
INTERNAL STANDARDS: TUNING:
OTHER:
SUMMARY: nothing found in samples despite high surrogate + Ms recoveries



VOLATILES SURROGATE RECOVERY

Primary Analysis Instrument ID: VAR 3600

	Date	Laboratory Reference No.	EPA-8020 & 602 BTEX Surrogate Recovery Fluorobenzene
01	03-20-91	Method Blank	128
102	03-21-91	Method Blank	126
03	03-21-91	29046 - 1	151
04	03-21-91	29046 - 2	125
05	03-21-91	29046 - 3	112
06	03-21-91	29046 - 4	285*
06	03-21-91	29046 - 4RE	279*
07	03-21-91	29046 - 5	130
80	03-21-91	29046 - 6	131
109	03-22-91	29046 - M04	177*
10	03-22-91	29046 - D04	232*

	1	Acceptane Range		
	Surrogate Standard	Water	Soil	
_	Fluorobenzene	60 - 132	53 - 152	

EPA-8080 & 602 Surrogate standards reported as percent recovery.

NA = Not Analyzed.

Comments: * Outside calibration range.

Approved By:

kdh.H

	·	The state of the s
CASE: 29046 SITE: 4	ANALYSIS: ρΑ	$H = \frac{1}{2} \mathbb{R}^{n}$
CASE: 29046 SITE: 4 LAB NOTES: indexs (123.cd) pyrene (123.cd) coelutes	w/ dibersp(a,h)enthraces	4
<u> </u>		
HOLDING TIMES: 3/13 ext - 3/7 se	mp. = ok	
NUMBER OF SAMPLES: Soil 3	Water _ /	
CALIBRATION Initial: sengo (q, h, i) perylene 30.	18 % RSD (confin = 7% RS)	رد
REF OR	·	
Continuing: beaco (9, h, i) perylene	- 40.6 %D (confin 32.6%.	b)
	· ·	52 A March
COEFFICIENT: not sue		
BLANKS Method:mo_lito		
Field: no hito		
FIELD DUPLICATES: not requested		
SURROGATES:		
MS/MSD: not requested		
INTERNAL STANDARDS:	Surrogate control Clark TUNING: No control charts	yet for soils
OTHER:	71	en wide
SUMMARY: no lits in samples		
		·
	-	
·		

SAMPLING REPORT FROM THE ALABAMA HIGHWAY DEPARTMENT



GOVERNOR

STATE OF ALABAMA HIGHWAY DEPARTMENT

MONTGOMERY, ALABAMA 36130

ROYCE G. KING HIGHWAY DIRECTOR

December 11, 1990

Mr. Gary Hinkle, Chief Installation Restoration Programs Branch Environmental Division National Guard Bureau Andrews Air Force Base, D.C. 20331-60080

Re: Project No. 51-006-003-008-901

US 80/Montgomery, AL

Dear Mr. Hinkle:

As you may recall, we wrote you on May 24, 1990, (copy attached) regarding highway construction activities along US 80 which is located adjacent to Dannelly Field ANGB. Enclosed you will find a copy of information generated as a result of our sampling activities. As you will note, we did not find that a major problem existed in this area with the exception that it appeared to us that the underground fuel storage next to the AASF hangar was leaking into the adjacent creek.

Per our May 24th letter, I am enclosing a copy of an invoice recently received from the ADEM Lab for sample analysis. Remittance for these costs should be submitted to the Alabama Highway Department at the above referenced address. Please reference the above project number and submit to my attention.

The invoices are misleading in that the total per sample is not in dollars as indicated, but is work time units. To convert these charges to dollars multiply these charges by 1.19. The total work time units is 4,732 which equates to \$5,631.08.

Should you have questions, please feel free to contact this Office.

Yours very truly,

Larry Lockett

Materials and Tests Engineer

BEC:sl

Attachment (3)

cc: File





STATE OF ALABAMA HIGHWAY DEPARTMENT

MONTGOMERY, ALABAMA 36130

May 24, 1990

ROYCE G. KING HIGHWAY DIRECTOR

Mr. Gary Hinkle, Chief Installation Restoration Programs Branch Environmental Division National Guard Bureau Andrews Air Force Base, D.C. 20331-6008

Dear Mr. Hinkle:

This letter is to confirm a telephone conversation of Monday, May 21, between Mr. David Hippensteel of your staff and Mr. B. E. Cox, a member of my staff. The topic discussed was potentially contaminated state right-of-way located along US 80 which is adjacent to Dannelly Field, Montgomery, Alabama.

During past operations it appears that various compounds may have drained from both the Air Guard and Army Guard Bases on Dannelly to State property. In order to determine if any contamination exist on State property from these facilities, sampling will be conducted in accordance with contract No. DE-AC05-840R21400 for sites 3 and 5. Additional sampling will be conducted along the perimeter fence in the vicinity of sites 3 and 5 where disposal may have occurred. The above referenced sites are associated with the Air Guard Facility.

In order to ensure that past operational and disposal practices by the Army Guard have not impacted State property, additional sampling will be conducted along the perimeter fence. Also, each major drainage ditch serving the Army Guard will be sampled.

As explained to Mr. Hippensteel it is imperative that this work be accomplished as soon as possible in order to prevent safety problems associated with highway construction in the area. Therefore, this Division will sample these areas in the near future. It is my understanding that NGB will reimburse the State for the cost of this work. Sample results will be forwarded to NGB for review. Should this sampling indicate an environmental problem, then further coordination relative to corrective action will be required.

Should you have questions, please feel free to contact this Office.

Yours very truly,

Fany Forbitt
Larry Wockett

Materials and Tests Engineer

BEC:s1

CC: Mr. Wm. J. Hartzog Mr. Mitch Kilpatrick



STATE OF ALABAMA HIGHWAY DEPARTMENT

MONTGOMERY. ALABAMA 36130

GUY HUNT GOVERNORE MORANDUM

ROYCE G. KING HIGHWAY DIRECTOR

DATE:

July 9, 1990

TO:

Larry Lockett

Materials & Test Engineer

FROM:

Bernard E. Cox, Jr. B& Environmental Engineer

SUBJECT: Project: 51-006-003-008-901

US 80 At Dannelly Field

On Friday, May 25, 1990, the writer collected 19 samples (soil and surface water), on right-of-way adjacent to Dannelly Field. The samples were collected to determine if past disposal practices by Army and Air Guard units at Dannelly had resulted in the creation of hazardous material sites on AHD ROW. This issue was raised by National Guard Bureau (NGB) in a letter to AHD dated May 3, 1990. In this letter NGB indicated that AHD would become responsible for any required remedial action should construction activities commence on potential sites.

Follow-up phone conversation and correspondence with NGB resulted in their agreeing to pay for sampling cost along the Dannelly-AHD ROW. Additionally, NGB was informed and now agrees that should remedial action be required, they (NGB) would be responsible for paying corrective action cost.

The writer was met at the sampling site by Johnny Day, project manager, for AHD. Sampling began at the west end of the area in question, on the west-side of Ft. Rufus Shepard. Four samples were collected between Station 423+48 and 420+85. No unusual traits were noted in this area. All soil samples were composed of numerous grab samples each 0 to 12 inches deep. These samples were composited in a metal pan lined with aluminum foil. Sampling spades were decontaminated with distilled water between sampling

Sampling progressed from Station 420+63 to Station 411+79. Seven samples were collected in this area. A noticeable kerosene odor was detected in the stream adjacent to the AASF hangar. Sample 2-3-4 was a water sample taken approximately 10 feet from the Guard frontage road in the drainage ditch. Samples 2-5-1 and 2-6-3/5 were also collected from this drainage course. Guard personnel informed the writer that a JP-4 UST located next to the drainage course had been emptied the morning of the 25th.

Sampling continued eastwardly from Station 411+59 to Station 401+60. Four samples were collected in this area with no unusual circumstances encountered. A composite sample from Station 411+59 to Station 404+91 was collected as was a composite sample from Sta. 404+61 to Sta. 404+60. Discreet samples were collected in the creek between Stations 401+60 and

Mr. Larry Lockett July 9, 1990 Page 2

Sampling progressed from Station 401+42 to Station 390+06, the east end point of the sampling area. Four samples were taken in this area with 4-1-1/3/5 being a composite sample along the old fence line. Three discreet samples were collected in the ditch outfall on the east end of the sampling zone.

The following numbering system for the samples was developed:

A-B-C/D/E.

A - Position signifies west to east with 1 on west end and 4 on the east end.

B- Sequential number, from 1 to 7. C/D/E - Position number with following values - (1- TPH); (2-BTEX-Soil); (3-Base Neutrals); (4-BTEX-Water); (5-Lead).

Sampling Protocol- Three types of samples were collected during the sampling Phase. These were composite soil, composite sludge and water. Composite soil samples were taken along the old fence line which is presently being moved by AHD contractors. Samples of this type were collected by taking a discreet sample every 10 to 15 feet along the fence and compositing by mixing in a lined pan. These samples were normally taken from 0 to 12 inches using aluminum scoops which were decontaminated with distilled water between sampling events. Samples collected in this manner were 1-2-1; 1-6-3/5; 2-1-1/3/5; 2-7-1/3/5; 3-1-1/3/5; 3-2-1/3/5; and 4-1-1/3/5.

Composite sludge soil samples were collected from four drainage courses exiting Dannelly onto AHD right-of-way. Either two or three discreet points were consolidated to make one sample. Materials were collected with an aluminum scoop and placed directly into a pint jar. Jars were filled to maximum capacity and sealed using aluminum foil and a screw on cap. Samples 1-3-1; 1-5-2; 2-2-1; 2-4-2; 2-5-1; 2-6-3/5; 3-3-2; 4-3-2; 4-3-3/5; and 4-4-1, were collected in this manner.

Two water samples, 2-3-4 and 3-4-4, were collected for volatile analysis. Volatile bottles secured from the ADEM laboratory were used for collection purposes. Only two of the four drainage courses exiting Dannelly had flow at the time of sampling. Samples were collected by placing bottles in the flow path and filling until no bubbles were left in the bottles. Bottle tops were then placed on the samples which were then checked to insure zero head space.

Samples for TPH analysis were taken to the AHD Environmental Lab and relinquished by the writer to Pat McCartha. Chain of custody was maintained and relinquished as required. All other samples were taken to the ADEM Lab and relinquished using appropriate chain of custody.

Sample results were received over a period of time from June 8 through July 1, 1990. Copies of the sample results are attached for your review. The following is a summary of the attached results:

June 5, 1990 - BTEX-ADEM Lab - Only sample 2-3-4 showed any volatile compounds. This sampled showed 469.7 PPB of 1-1-1- Trichlorethane. This

Mr. Larry Lockett July 9, 1990 Page 3

sample was the water sample in the creek next to the AASF where kerosene was noted at the time of sampling. The completed BTEX-Volatile sampling was forwarded June 8, 1990, from ADEM to this Office. Only sample 2-4-2, a composite soil sample in the same creek, showed any compounds. This sample showed .19 ug\l'of 1,1,1, - Trichloroethane.

June 8, 1990 - TPH results -AHD Lab - Eight samples were analyzed for TPH with all eight showing some signs of TPH. This was to be expected since the samples were taken along the fence line bordering a major highway where diesel fuel and other fuels had been used by the Guard for weed control. The highest TPH value (575 ppm) was a composite sample taken between Stations 413+90 and 411+79. This is the fence line immediately in front of the AASF Hangar. The next highest reading was found at Station 390+83. This sample was collected next to a sorbent boom in the drainage ditch on Guard property.

June 19/26, 1990 - Base Neutrals; Lead; EPTOX Metals - No regulated EPT metals (including lead) were found in any samples. All total lead levels were low and consistent with normal soil conditions. No base neutrals were observed in any samples other than 4-4-3/5. Low levels (PPB) of various base neutral compounds were found in this sample which was collected in the drainage ditch on the east end of the project.

<u>Conclusions</u> - Diesel fuel was known to be used as a weed killer by Guard personnel along the fence line so finding TPH in this area was expected. No base neutrals, metals or volatiles were found in these samples. Thus, none of this area would constitute a hazardous waste or other regulated site. It should be noted that TPH is an indicator only and has regulatory value only when used in conjunction with an UST.

Only two questionable areas were noted and they are both drainage ditches exiting Guard property onto AHD ROW. These areas are adjacent to UST operated by Guard units. TPH and some indicators parameters, at very low levels, were found in two samples. Neither site should impact AHD construction activities.

Recommendations - Sampling results should be forwarded to NGB. Construction activities should continue as no regulated sites were noted on AHD ROW.

BEC:s1

cc: Mr. Mitch Kilpatrick

File



STATE OF ALABAMA HIGHWAY DEPARTMENT

MONTGOMERY, ALABAMA 36130

June 8,1990

ROYCE G. KING HIGHWAY DIRECTOR

MEMORANDUM

TO:

Mr. Stanley R. Armstrong

Assistant Materials & Tests Engineer

FROM:

Pat McCartha

RE:

Test results for Petroleum Hydrocarbon

Contamination; Project 51-006-003-901 Montgomery

county

Attached are test results for samples taken from various locations at the referenced site. Copies are being sent to the Division for their information.

wpm

cc: Division

Mr. Bill Holmes

File

ALABAMA HIGHWAY DEPARTMENT

BMT-16 Rev.

COPIES TO. Chemical Lab	PROJECT NO(S). 51-006-003-901
••••••	COUNTY Montgomery
	DIVISION 6
••••••	DATE June 8, 1990
1 TUODECHE CO.	

) INSPECTION

REPORT OF)ANALYSIS ON SAMPLE OF.. Soil for Total Petroleum Hydrocarbon

SOURCE OF MATERIAL.... See Station & Offset SAMPLED BY / DATE.... B. Cox; 5/25/90 SUBMITTED BY / DATE... B. Cox; 05/25/90 Received By/ Date.... Strickland; 05/25/90 Tested By/ Date... D. Hicks; 06/06/90 ADEM Reference No.... REMARKS...

TEST RESULTS

		IESI KESULIS	
LAB NO.	MARKS	MATERIAL	TOTAL PETROLEUM HYDROCARBON
3q1-1918 3q1-1919 3q1-1920 3q1-1921 3q1-1922 3q1-1923 3q1-1924 3q1-1925	1-2-1 2-1-1 2-5-1 2-7-1 3-1-1 3-2-1 4-1-1 4-4-1	Soil Soil Soil Soil Soil Soil Soil	24 77 159 575 207 82 44 456

Total Petroleum Hydrocarbon results in mg/kg on a dry weight basis

ADEM

ALABAMA DEPARTMENT OF ENVIRONMENTAL MANAGEMENT

Attached are laboratory results from samples submitted to the ADEM

Central Laboratory for analysis by the Highway Department.



Guy Hunt Governor

Leigh Pegues, Director

1751 Cong. W. L.

Dickinson Drive June 26, 1990

Montgomery, AL 36130

205/271-7700

MEMORANDUM

TO:

FROM:

SUBJECT:

Buddy Cox

Field Offices:

Highway Department

Laboratory Results

Unit 806, Building 8

225 Oxmoor Circle

225 Oxmoor Circle Birmingham, AL

35209

205/942-6168

205/942-6168

P.O. Box 953 Decatur, AL

35602

205/353-1713

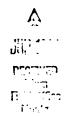
JC/mpt

Attachments

2204 Perimeter Road Mobile, AL 36615 205/479-2336 .

- SAMPLE ANALYSIS REPORT - 06/22/90

To: ALABAMA HIGHWAY DEPARTMENT



Attn: BUDDY COX

Lab number : 0105303 / Sample number : HIWAY

Sample matrix : SOIL

Report Date: 06/22/90

COLLECTION INFORMATION

Date/Time/By: 05/25/90 9:30 COX Location : DANNELLY ANGB, 1-6-3/5

ADEM CENTRAL LABORATORY - RESULTS REPORT -

June 22, 1990

			•	
Lab#	Test	Result	UnitsDL*	Analdate
Lab# 0105303	1,2,4Trichlorobenzene 1,2-Dichlorobenzene 1,2-Diphenylhydrazine 1,3-Dichlorobenzene 1,4-Dichlorobenzene 2,3,7.8-Tetrachlorodibe 2,4-Dinitrotoluene 2,6-Dinitrotoluene 2-Chloronaphthalene 3,3'-Dichlorobenzidine 4-Bromophenyl phenyl eh 4-Chlorophenyl phenyl eh Acenaphthalene Silver-EP Anthreene Arsenic-EP Benzo(a)anthacene Barium-EP	0.33 0.33 0.33 0.33 0.33 0.33 0.33 0.33	ug/L U	05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90
	Benzo(a)pyrene Benzo(b)fluoanthene Butyl benzyl phthalate	0.33 0.33	mg/L U ug/L U ug/L U ug/L U	06/13/90 05/28/90 05/28/90 05/28/90

U denotes results less than the instrument detection limit.

STATE OF ALABAMA DEPARTMENT OF ENVIRONMENTAL MANAGEMENT MONTGOMERY, ALABAMA

LABORATORY:	(X Mont	gomery	() Mobile	() Birmi	ngham
<u> </u>	Potable Water [] Surface Water [] Soil/Sediment [] Wastewater []	Hazardous Wast Groundwater	esite [] Ignita	ivity []	Composite Grab Container P G
ource <u>Par</u>	welly ANGB	US80			/
Location	-6-3/5				
) Discharge	from		to		
_	(Po	oint Source)		(Receiving	
omments			Preservative(s)	<u> </u>	
ьн	D.O Sp.	CondPARAM	Salinity	Turb.	
Date (mg/1)		Value (mg/l)	Date Val (mg/l)		Value mg/l)
Acid		Phenol	Al		Mn
ALK		PO, -P	Ag		Na
BOD ₅ -		(s=)			Ni
(C1 ⁻)_		(so ₄ =)	Ba		Pb
COD _		TSS	Ca		Pt
CN		TDS	Cd		Sb
(F ⁻)_		TFS	Cr ^T		Se
Hard _		TKN	Cr ⁺⁶		Zn
NH ₃ -N _		TOC	Cu	Other	
NO ₃ -N _		TON	Fe	Base 14	entrels
NO ₂ -N _		TS	Нд	head (Tof 4 EPT)
0 & G _	·	vss	Mg		
_	ſ			F. Coli.	
B.Cox	3.6/.	9:30 \$125/90	B. Cox	3 5///	2:00 5/25/90
AMPLE COLLECT	ED BY (Signature	e) DATE/TIME	RELINQUISHED BY	(Signature)	DATE/TIME /
ECEIVED BY	(Signature	DATE/TIME	RELINQUISHED BY	(Signature)	DATE/TIME
ECEIVED BY, ECEIVED IN LA	Signature Show 5/25 B BY (Signature	2:00	RELINQUISHED BY LABORATORY I.D.	(Signature) 0105303 NO.	DATE/TIME
SEND REPORT TO	α	Lovey Mato	Test)		<u> </u>
ADEM Form 68 5	5/83		tal Non-Filtrable Resi	due IS	= Total Residue

ADEM Form 68 5/83

ADEM CENTRAL LABORATORY

- SAMPLE ANALYSIS REPORT - 06/01/90

To: ALABAMA HIGHWAY DEPARTMENT



Attn: BUDDY COX

Lab number : 0105307 Sample number : HIWAY

Sample matrix : SOIL

Report Date: 06/01/9

COLLECTION INFORMATION

Date/Time/By: 05/25/90 9:10 COX Location : DANNELLY ANGB US80, 1-5-2

ADEM CENTRAL LABORATORY
- RESULTS REPORT -

June 1, 1990

Lab#	Test	Result	UnitsDL*	Analdate
	1,1,1,2-Tetrachloroetha 1,1,1-Trichloroethane 1,1,2,2-Tetrachloroetha 1,1,2Trichloroethane 1,1-Dichloroethane 1,1-Dichloroethylene 1,1-Dichloropropene 1,2,3-Trichlorobenzene 1,2,3-Trichloropropane 1,2,4-Trichlorobenzene 1,2-Dichloropropane 1,2-Dichloropropane 1,3-Trimethylbenzene 1,3-Dichloropropane	0.0500 0.0500 0.0500 0.0500 0.0500 0.0500 0.0500 0.0500 0.0500	ug/g U	05/31/90 05/31/90

U denotes results less than the instrument detection limit.

Lab#	Test	Result	UnitsDL*	Analdate
0105307	Bromomethane	0.0500	ug/g U	05/31/90
	cis-1,2-Dichloroethylen		ug/g U	05/31/90
	Chlorobenzene	0.0500	ug/g U	05/31/90
	Chlorodibromomethane	0.0500	ug/g U	05/31/90
	Chloroethane	0.0500	ug/g U	05/31/90
	Bromoform	0.0500	ug/g U	05/31/90
	Chloroform	0.0500	ug/g U	05/31/90
	Chloromethane	0.0500	ug/g U	05/31/90
	Carbon Tetrachloride		ug/g U	05/31/90
	Dibromomethane	0.0500	ug/g U	05/31/90
	Dichlorofluoromethane	0.0500	ug/g U	05/31/90
	Dichloromethane	0.0500	ug/g U	05/31/90
	Ethylbenzene	0.0500	ug/g U	05/31/90
	Fluorotrichloromethane	0.0500	ug/g U	05/31/90
	Hexachlorobenzene		ug/g U	05/31/90
	Isopropylbenzene	0.0500	ug/g U	05/31/90
	m-Dichlorobenzene		ug/g U	05/31/90
	m-Xylene	0.0500	ug/g U	05/31/90
	Naphthalene		ug/g U	05/31/90
	n-Butylbenzene		ug/g U	05/31/90
	n-Propylbenzene		ug/g U	05/31/90
	o-Chlorotoluene		ug/g U	05/31/90
	o-Dichlorobenzene		ug/g U	05/31/90
	o-Xylene		ug/g U	05/31/90
	p-Chlorotoluene		ug/g []	05/31/90
•	p-Dichlorobenzene		ug/g U	05/31/90
		0.0500		05/31/90
	p-Xylene		ug/g U	05/31/90
	Secbutylbenzene		ug/g U	05/31/90
	Styrene	0.0500		05/31/90
	t-1.2Dichloroethane		ug/g U	05/31/90
	Tertbutylbenzene		ug/g U	05/31/90
	Trichloroethylene	0.0500		05/31/90
	Toluene		ug/g U	05/31/90
	Vinyl Chloride	0.0500	ug/g U	05/31/90

U denotes results less than the instrument detection limit.

STATE OF ALABAMA DEPARTMENT OF ENVIRONMENTAL MANAGEMENT MONTGOMERY, ALABAMA

LABORATORY:	✓ Montg	gomery	() Mobile	() Bir	mingham
Sample Type:	Potable Water [] Surface Water [] Soil/Sediment [] Wastewater []	Landfill Leach Hazardous Wast Groundwater Waste (Special Ha	esite [] Igni [] Corr	c Extraction [tability [cosivity [tivity [Composit 2) Grab Container P G
Source	Dannelly Al	6B 4580)		,
Location /-	5-2				4
() Discharge	from		to		
Comments		int Source)	_	(Receivin	g Water)
			_ Preservative(s		
pH	D.O Sp.	CondPARAM	Salinity METERS	Turb	
Date	Value Date			alue Date	Value
(mg/1)	(mg/1)	(mg/1)	arde Date	(mg/l)
Acid -	P	henol	Al		Mn
ALK _	P	O ₄ -P	Ag		Na
BOD ₅ -		(s ⁼)	As		Ni
(C1 ⁻)	(so ₄ =)	Ba		Pb
COD		rss	Ca		Pt
CN		TDS	Cd		Sb
(F ⁻)_		TFS	Cr ^T		Se
Hard _		TKN	Cr ⁺⁶		Zn
NH ₃ -N -		гос	Cu	Other	
NO ₃ -N _		ron	Fe		(1/1/
NO ₂ -N _		TS	Нд		
0 & G _		vss	Mg		
	• .	•		F. Coli	
B 66	D. GX	9:10 5/25/90	7361	3. Cv	2:00 5/25/91
SAMPLE COLLECT	ED BY (Signature)	9:10 5/25/90 DATE/TIME	RELINQUISHED BY	(Signature)	DATE/TIME
	•				
RECEIVED BY	(Signature)	DATE/TIME	RELINQUISHED BY	(Signature)	DATE/TIME
RECEIVED BY RECEIVED IN LA	Signature) B BY (Signature)	25 2:00	RELINQUISHED BY	110530	DATE/TIME
SEND REPORT TO	· Cox (Me)	+ Test)			

TSS = Total Non-Filtrable Residue
TOS = Total Filtrable Residue

TS = Total Residue
VSS = Volatile Residue

ADEM CENTRAL LABORATORY

- SAMPLE ANALYSIS REPORT - 05/31/90

To: ALABAMA HIGHWAY DEPARTMENT

The state of the s

Attn: BUDDY COX

Lab number : 0105297 Sample number : HIWAY

Sample matrix : SOIL

COLLECTION INFORMATION

Date/Time/By: 05/25/90 9:20 COX Location : DANNELLY ANGB, 1-3-2

ADEM CENTRAL LABORATORY

- RESULTS REPORT -

May 31, 1990

Report Date: 05/31/9

Lab#	Test	Result UnitsDL* Analdate
0105297	1,1,1,2-Tetrachloroetha 1,1,1-Trichloroethane	0.0500 ug/g U 05/29/90 0.0500 ug/g U 05/29/90
-	1,1,2,2-Tetrachloroetha	0.0500 ug/g U 05/29/90
	1.1,2Trichloroethane	0.0500 ug/g U 05/29/90
	1,1-Dichloroethane	0.0500 ug/g U 05/29/90
	1,1-Dichloroethylene	0.0500 ug/g U 05/29/90
	1,1-Dichloropropene	0.0500 ug/g U 05/29/90
	1,2,3-Trichlorobenzene	0.0500 ug/g U 05/29/90
	1,2,3-Trichloropropane	0.0500 ug/g U 05/29/90
	1,2,4-Trichlorobenzene	0.0500 ug/g U 05/29/90
	1,2,4-Trimethylbenzene	0.0500 ug/g U 05/29/90
	1,2-Dicholoethane	0.0500 ug/g U 05/29/90
	1.2-Dichloropropane	0.0500 ug/g U 05/29/90
	1.3.5-Trimethylbenzene	0.0500 ug/g U 05/29/90
	1.3-Dichloropropane	0.0500 ug/g U 05/29/90
	1,3-Dichloropropene	0.0500 ug/g U 05/29/90
-	2,2-Dichloropropane	0.0500 ug/g U 05/29/90
	Tetrachloroethylene	0.0500 ug/g U 05/29/90
	Bromobenzene	0.0500 ug/g U 05/29/90
	Bromochloromethane	0.0500 ug/g U 05/29/90
	Bromodichloromethane	0.0500 ug/g U 05/29/90
	Benzene	0.0500 ug/g U 05/29/90

^{*} U denotes results less than the instrument detection limit.

0105297 Bromomethane 0.0500 ug/g cis-1,2-Dichloroethylen 0.0500 ug/g Chlorobenzene 0.0500 ug/g Chlorodibromomethane 0.0500 ug/g Chloroethane 0.0500 ug/g Bromoform 0.0500 ug/g Chloroform 0.0500 ug/g Chloromethane 0.0500 ug/g Carbon Tetrachloride 0.0500 ug/g	:sDL*	Analdate
Dibromomethane Dichlorofluoromethane Dichlorofluoromethane Dichloromethane Dichloromethane Dichloromethane Ethylbenzene Fluorotrichloromethane Hexachlorobenzene Isopropylbenzene M-Dichlorobenzene M-Dichlorobenzene M-Zylene M-Butylbenzene M-Propylbenzene M-Sylene M-P-Chlorotoluene M-Dichlorobenzene M-Dichlorobenzene M-Isopropyltoluene M-Isopropyltol		05/29/90 05/29/90

^{*} U denotes results less than the instrument detection limit.

	non (G	UMERY, ALABAMA	
LABORATORY:	Montgomery	() Mobile	() p;
Sample Type:	Potable Water [] Landfill Lea		() Birmingham
	Surface Water [] Hazardous Wa	Astesire [] Tame	traction [] Composite
1	Soil/Sediment [X] Groundwater Wastewater [] Waste (Special	[] Corrosiv	ity [] Contains
Source Da	innelly ANGB USBO	Handling) [] Reactivi	ty [] G
	- 3-2		
() Discharge f	rom		
	(Point Source)	to	
Comments		Preservative(s)	(Receiving Water)
рН D	.0 Sp. Cond	Solini.	
	FAR	SalinityTur	b
Date Va (mg/1)	alue Date Value	Date Value	D.
Acid	(mg/1)	(mg/l)	Date Value (mg/l)
ALK	Pheno I		Mn
BOD ₅	PO ₄ -P	Ag	Na Na
(C1 ⁻)	(S ⁼)	As	Ni
COD	(SO ₄ =)	Ba	Pb
CN	TSS	Ca	Pt
(F ⁻)	TES	Cd	Sb
Hard	TFS	Cr ^T	Se
NH ₃ -N	TKN TOC	Cr ⁺⁶	Zn
NO3-N	TOC	Cu	Other
NO ₂ -N		Fe	BIEX-VO
0 & G	TS VSS	Нд	
	vss	Mg	
7 6x	71000	~ . 4	F. Coli
SAMPLE COLLECTED	B. 64 9:20 5/25/50 BY (Signature) DATE/TIME	B. Cox B.	ex 2:00 5/25-190
	, Date, Time	RELINQUISHED BY (Si	gnature) DATE/TIME
RECEIVED BY	(Signature) DATE/TIME	PEL INQUIRCUM	
	- July IIME	RELINQUISHED BY (Si	gnature) DATE/TIME
RECEIVED BY	(Signature) DATE/TIME	RELINQUISHED BY (Si	· · · · · · · · · · · · · · · · · · ·
X Ollino	15/25 21/2		gnature) DATE/TIME
RECEIVED IN LAB BY	(Signature) DATE/TIME	LABORATORY I.D. NO.	5297
END REPORT TO:	Cox (Mot & Test	2)	
DEM Form 68 5/83	TSS - Tot	al Non-Filtrable Residue	70
30 3/03	TDS = Tot	al Filtrable Residue	TS = Total Residue

LABORATORY:	Montgomery	() Mobile	() Birmingham
Sample Type:	Soil/Sediment Groundwater	esite [] Ignita	Extraction Composite bility [] Grab
Source Do	unelly ANGB US80		1
Location 2	-7-3/5		
() Discharge	from (Point Source)	to	
Comments	(Point Source)	_ Preservative(s)	(Receiving Water)
pH	D.O Sp. Cond		
	Value Date Value	Date Val	ue Date Value
(mg/l)	\g, -,	(mg/1)	(mg/1)
Acid _	Phenol	A1	Mn
ALK _	PO ₄ _P	Ag	Na
BOD_5 -	(s ⁼)	As	Ni
(C1 ⁻)_	(so ₄ =)	Ba	
COD _	TSS	Ca	Pt
CN	TDS	Cd	Sb
(F ⁻)_	TFS	Cr ^T	Se
Hard _	TKN	Cr ⁺⁶	Zn
NH ₃ -N _	тос	Cu	Other
NO ₃ -N _	TON	Fe	Base Newforls
NO ₂ -N _	TS	Нд	Pb (Tut + CPT)
O & G _	vss	Mg	
B. Lox SAMPLE COLLECT	D. E. (//: 1/25/50 ED BY (Signature) DATE/TIME	B. E RELINOVISHED BY	F. Coli. B. E. 2:0007/25795 (Signature) DATE/TIME
RECEIVED BY	(Signature) DATE/TIME	RELINQUISHED BY	(Signature) DATE/TIME
RECEIVED BY RECEIVED IN LA		RELINQUISHED BY LABORATORY I.D. 1	(Signature) DATE/TIME
SEND REPORT TO	: _ Cox (Mat + Test)		

- SAMPLE ANALYSIS REPORT - 06/13/90

To: ALABAMA HIGHWAY DEPARTMENT

Attn: BUDDY COX

Lab number : 0105299 Sample number : HIWAY

Sample matrix : SOIL

Report Date: 06/13/90

COLLECTION INFORMATION

Date/Time/By: 05/25/90 11:10 COX Location : DANNELLY ANGB, 2-7-3/5

ADEM CENTRAL LABORATORY
- RESULTS REPORT -

June 13, 1990

Lab#	Test	Result	UnitsDL*	Analdate
0105299	1,2,4,-Trichlorobenzene 1,2-Dichlorobenzene 1,2-Diphenylhydrazine 1,3-Dichlorobenzene 1,4-Dichlorobenzene 2,3,7,8-Tetrachlorodibe 2,4-Dinitrotoluene 2,6-Dinitrotoluene 2-Chloronaphthalene 3,3'-Dichlorobenzidine 4-Bromophenyl phenyl eh 4-Chlorophenyl phenyl e Acenaphthalene Acenaphthalene Acenaphthalene Acenaphthalene Acenaphthalene Acenaphthalene Bariue-EP Benzo(a)anthacene Barium-EP Benzo(a)pyrene	0.33 0.33 0.33 0.33 0.33 0.33 0.33 0.33	UnitsDL* ug/L U Analdate 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 06/11/90 05/28/90 06/11/90 05/28/90 05/28/90	
	Benzo(b)fluoanthene Butyl benzyl phthalate	0.33	ug/L U ug/L U	05/28/90 05/28/90

^{*} U denotes results less than the instrument detection limit.

- SAMPLE ANALYSIS REPORT - 06/13/90

To: ALABAMA HIGHWAY DEPARTMENT

JUH 1990 RECEIVED Adem Field Office Montg.

Attn: BUDDY COX

Lab number : 0105300 Sample number : HIWAY Sample matrix : SOIL

Report Date: 06/13/90

COLLECTION INFORMATION

Date/Time/By: 05/25/90 11:28 COX Location : DANNELLY ANGB, 2-6-3/5

ADEM CENTRAL LABORATORY
- RESULTS REPORT -

June 13, 1990

Lab#	Test	Result	UnitsDL*	Analdate
0105300	1.2,4,-Trichlorobenzene 1.2-Dichlorobenzene 1,2-Diphenylhydrazine 1,3-Dichlorobenzene 1,4-Dichlorobenzene 2,3,7.8-Tetrachlorodibe 2,4-Dinitrotoluene 2,6-Dinitrotoluene 2-Chloronaphthalene 3,3'-Dichlorobenzidine 4-Bromophenyl phenyl eh 4-Chlorophenyl phenyl eh 4-Chlorophenyl phenyl e Acenaphthalene Acenaphthalene Silver-EP Anthroene Arsenic-EP Benzo(a)anthacene Barium-EP Benzo(b)fluoanthene Butyl benzyl phthalate	0.33 0.33 0.33 0.33 0.33 0.33 0.33 0.33	ug/L U	05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 06/11/90 05/28/90 06/11/90 05/28/90 06/11/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90

^{*} U denotes results less than the instrument detection limit.

Lab#	Test	Result	UnitsDL*	Analdate
0105300	Bis (2-chlororthyl) eth	0.33	ug/L U	05/28/90
	Bis(2-chloroethoxy)meth	0.33	ug/L U	05/28/90
	Bis (2-Chloroisopropyl)	0.33	ug/L []	05/28/90
	Bis(2-ethylhexyl)phthal		ug/L U	05/28/90
	Benzo(g,h,i)perylene		ug/L U	05/28/90
	Benzidine		ug/L U	05/28/90
	Benzo(k)fluoranthene		ug/L U	05/28/90
	Cadmium-EP		mg/L U	06/11/90
	Chromium-EP		mg/L U	06/11/90
	Chrysene		ug/L U	05/28/90
	Dibenzo(a.h)anthracene		ug/L U	05/28/90
	Dibutyl phthalate		ug/L U	05/28/90
	Diethyl phthalate		ug/L U	05/28/90
	Dimethylphthalate		ug/L U	05/28/90
	Di-n-octyl phthalate		ug/L U	05/28/90
	Fluoranthene		ug/L U	05/28/90
•	Fluorene		ug/L U	05/28/90
*	Hexachlorobutadiene		ug/L U	05/28/90
	Hexachlorobenzene		ug/L U	05/28/90
	Hexachlorocyclopentadie	0.33	IJ	05/28/90
	Hexachloroethane		ug/L U	05/28/90
*	Mercury-EP		mg/L U	06/11/90
	Isophrone		ug/L U	05/28/90
	Indeno(1,2,3-cd)pyrene		ug/L U	05/28/90
	Naphthalene		ug/L U	05/28/90
	Nitrobenzene		ug/L U	05/28/90
	N-nitroso-di-n-propylam		ug/L U	05/28/90
	N-nitrosodimethylamine		ug/L U	05/28/90
	N-nitrosodiphenylamine		ug/L U	05/28/90
•	Pyrene		ug/L U	05/28/90
•	Phenanthrene		ug/L U	05/28/90
	Lead-EP		mg/L U	06/11/90
	Lead in Soil		ug/g U	06/11/90
•	Selenium-EP	0.01	mg/L U	06/11/90

^{*} U denotes results less than the instrument detection limit.

Hi-way

STATE OF ALABAMA DEPARTMENT OF ENVIRONMENTAL MANAGEMENT MONIGOMERY, ALABAMA

Source	Dannelly ANGB USB	Wastesite [] Ignitabir [] Corrosival Handling [] Reactivi	ity [] Co
	2-6-3/5-		
() Discharge			
Comments	(Point Source)	to	(Pagaini in
		Preservative(s)	(Receiving Water)
	Sp. Cond.	Salinity	L
Date	Value non	RAMETERS	D
(mg/1)	Value Date Value (mg/l)	Date Value	Date Value
Acid	Pheno1	(mg/1)	Date Value (mg/1)
ALK	PO,P	A1	Mn
BOD_5	(s=)	Ag	Na
(C1 ⁻)	(SO ₄ =)	As	Ni
COD	TSS	Ba	Pb
CN	TDS	Ca	Pt
(F¯)	TFS	Cr ^I	Sb
Hard	TKN	Cr ⁺⁶	Se
NH ₃ -N	тос	Cu	Zn
NON	TON	Fe	Other
NO ₂ -N O & G	TS	Нg	Base Newfreb
	vss	Mg	Ph (TotIEPT)
B. COLLECTED	B. I. (Signature) DATE/TIME	RELINQUISHED BY (SE	F. Coli
CEIVED BY	(Signature) DATE/TIME	RELINQUISHED BY (Sig	nature) DATE/TIME
EIVED BY EIVED IN LAB BY D REPORT TO:	(Signature) DATE/TIME Ver 5/25 2:00 (Signature) DATE/TIME Cox Mod + Test	RELINQUISHED BY (Sig	nature) DATE/TIME 5300

The Chrometogram of this sample reneals a pattern of hydrocarbons that matches that found in kerosene.

- SAMPLE ANALYSIS REPORT - 06/13/90

To: ALABAMA HIGHWAY DEPARTMENT



Attn: BUDDY COX

Lab number : 0105302 Sample number : HIWAY

Sample matrix : SOIL

Report Date: 06/13/90

COLLECTION INFORMATION

Date/Time/By: 05/25/90 9:35 COX Location : DANNELLY ANGB, 2-1-3/5

ADEM CENTRAL LABORATORY
- RESULTS REPORT -

June 13, 1990

Lab#	Test	Result	UnitsDL*	Analdate
0105302	1,2,4,-Trichlorobenzene 1,2-Dichlorobenzene 1,2-Diphenylhydrazine 1,3-Dichlorobenzene 1,4-Dichlorobenzene 2,3,7,8-Tetrachlorodibe 2,4-Dinitrotoluene 2,6-Dinitrotoluene 2-Chloronaphthalene 3,3'-Dichlorobenzidine 4-Bromophenyl phenyl eh 4-Chlorophenyl phenyl eh 4-Chlorophenyl phenyl e Acenaphthalene Acenaphthalene Silver-EP Anthroene Arsenic-EP Benzo(a)anthacene Barium-EP Benzo(b)fluoanthene	0.33 0.33 0.33 0.33 0.33 0.33 0.33 0.33	ug/L U	05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 06/11/90 05/28/90 06/11/90 05/28/90 06/11/90 05/28/90 05/28/90 05/28/90 05/28/90
	Butyl benzyl phthalate	0.33	ug/L U	05/28/90

^{*} U denotes results less than the instrument detection limit.

LABORATORY:	Мо	ntgomery	() Mobile	() Bi	rmingham
Sample Type:	Potable Water [Surface Water [Soil/Sediment [Wastewater [j⊓Hazardous Was X Groundwater	tesite [] Ignit	sivity [Composite Grab Container P G
Source	Dannelly An	LB US80			
Location \mathcal{Z}					
() Discharge	from		to		
Comments		(Point Source)	Preservative(s)	(Receivi	
рН	D.OS		_ Salinity AMETERS	_ Turb	
(mg/1)		e Value (mg/1)	Date Va		e Value (mg/1)
Acid ALK		Phenol PO ₄ -P	A1		Mn Na
	-	(S ⁼)	Ag As		Ni Ni
(C1 ⁻)		(so ₄ =)			Pb
COD		TSS	Ca		Pt
CN-		TDS	Cd		Sb
(F ⁻)		TFS	- Cr ^T		Se
Hard		TKN	Cr ⁺⁶		Zn
NH ₃ -N		TOC	Cu	Other	
NO ₃ -N	·	TON	Fe	— Bara	· Neufrels
NO ₂ -N		TS	Нg	Pb	(Tot + EPT
0 & G	· · ·	VSS	Mg		
	_			F. Co	li
B.Cx SAMPLE COLLEC	B & (Signat	9:35 5/25/50 ure) DATE/TIME	RELINQUISHED BY	B 4 Y (Signature	2:00 35/25-/90 E) DATE/TIME
RECEIVED BY	(Signat	ure) DATE/TIME	RELINQUISHED BY	Y (Signature	DATE/TIME
RECEIVED BY	Signature (Signature)	ture) DATE/TIME	RELINQUISHED BY	Y (Signature	
RECEIVED IN I			LABORATORY I.D	. NO.	
South VELOKI I	··				

ADEM CENTRAL LABORATORY

- SAMPLE ANALYSIS REPORT - 06/01/90

To: ALABAMA HIGHWAY DEPARTMENT



Report Date: 06/01/9

Attn: BUDDY COX

Lab number : 0105298 Sample number : HIWAY

Sample matrix : SOIL

COLLECTION INFORMATION

Date/Time/By: 05/25/90 10:45 COX Location : DANNELLY ANGB, 2-4-2

ADEM CENTRAL LABORATORY
- RESULTS REPORT -

June 1, 1990

Lab#	Test	Result	UnitsDL*	Analdate
0105298	1.1,1.2-Tetrachloroetha 1.1.1-Trichloroethane 1.1.2.2-Tetrachloroetha 1.1.2Trichloroethane 1.1-Dichloroethane 1.1-Dichloroethylene 1.1-Dichloropropene 1.2.3-Trichlorobenzene 1.2.3-Trichloropropane 1.2.4-Trichlorobenzene 1.2.4-Trimethylbenzene 1.2-Dichloropropane 1.3-Dichloropropane	0.1900 0.0500 0.0500 0.0500 0.0500 0.0500 0.0500 0.0500 0.0500 0.0500 0.0500 0.0500 0.0500 0.0500 0.0500	ug/g U	05/31/90 05/31/90

^{*} U denotes results less than the instrument detection limit.

Lab#	Test	Result	UnitsDL*	Analdate
0105298	Bromomethane		ug/g U	05/31/90
	cis-1,2-Dichloroethylen		ug/g U	05/31/90
	Chlorobenzene		ug/g U	05/31/90
	Chlorodibromomethane		ug/g U	05/31/90
	Chloroethane		ug/g U	05/31/90
	Bromoform		ug/g U	05/31/90
	Chloroform		ug/g U	05/31/90
	Chloromethane		ug/g U	05/31/90
	Carbon Tetrachloride		ug/g U	05/31/90
	Dibromomethane	0.0500	ug/g U	05/31/90
	Dichlorofluoromethane	0.0500	ug/g U	05/31/90
•	Dichloromethane	0.0500	ug/g []	05/31/90
	Ethylbenzene	0.0500	ug/g U	05/31/90
	Fluorotrichloromethane	0.0500	ug/g U	05/31/90
	Hexachlorobenzene	0.0500	ug/g U	05/31/90
	Isopropylbenzene	0.0500	ug/g U	05/31/90
	m-Dichlorobenzene	0.0500	ug/g U	05/31/90
	m-Xylene	0.0500	ug/g U	05/31/90
	Naphthalene		ug/g U	05/31/90
	n-Butylbenzene		ug/g U	05/31/90
	n-Propylbenzene	0.0500	ug/g U	05/31/90
	o-Chlorotoluene		ug/g U	05/31/90
	o-Dichlorobenzene	0.0500	ug/g U	05/31/90
•	o-Xylene	0.0500	ug/g U	05/31/90
•	p-Chlorotoluene	0.0500	ug/g U	05/31/90
	p-Dichlorobenzene	0.0500	ug/g U	05/31/90
	p-Isopropyltoluene	0.0500	ug/g U	05/31/90
	p-Xylene '	0.0500	ug/g U	05/31/90
	Secbutylbenzene	0.0500	ug/g U	05/31/90
	Styrene	0.0500	ug/g U	05/31/90
	t-1.2Dichloroethane	0.0500	ug/g U	05/31/90
	Tertbutylbenzene	0.0500	ug/g U	05/31/90
	Trichloroethylene	0.0500	ug/g U	05/31/90
	Toluene	0.0500	ug/g U	05/31/90
	Vinyl Chloride	0.0500	ug/g U	05/31/90

st U denotes results less than the instrument detection limit.

LABORATORY:	Montgoi	nery	() Mobile	() Birmingha	m
Sample Type:	Surface Water [] Soil/Sediment []	Landfill Leachat Hazardous Wastes Groundwater Waste (Special Hand	ite [] Ig [] Co	xic Extract nitability rrosivity activity	[] Gra	posite b tainer P G
Source	Dannelly ANGI	y 1580				
Location 2	. 1					
() Discharge			to			
•		nt Source)			ceiving Wate	r)
Comments						
рН	D.O Sp. C	ondSa PARAME		Turb		
	Value Date	Value	Date	Value	Date	Value
(mg/1)		g/1)	(mg/1)		(mg/1	
Acid		enol	A1		Mn	
ALK	~	_P	Ag		Na	
BOD_5 -		s ⁼)	As		Ni	
(C1_)	(SC) ₄ =)	Ba		Pb	
COD	T:	SS	Ca		Pt	
CN	TI	os	Cd		Sb	
(F¯)_	T	FS	Cr ^T		Se	
Hard	TI	(N	Cr ⁺⁶		Zn	
NH ₃ -N -	TO	oc	Cu		Other	
NO ₃ -N _	TO	ON	Fe		BIEX V	5/
NO2-N		rs	Hg			
O & G	V:	SS	Mg			
7.64	B. L. 41	10:45 5/25/90	B. E. 41	B. E	F. Coli	06 5/25-/9
SAMPLE COLLECT	TED BY (Signature)	DATE/ŤIMÉ	RELINQUZSHÆD	BY (Sign	ature) DAT	E/TIME
RECEIVED BY	(Signature)	DATE/TIME	RELINQUISHED	BY (Sign	nature) DAT	E/TIME
RECEIVED BY	(Signature)	DATE/TIME	RELINQUISHED		ature) DAT	E/TIME
RECEIVED IN LA	\sim \sim \sim \sim \sim	DATE/TIME	LABORATORY I	.D. NO.		

- SAMPLE ANALYSIS REPORT - 05/31/90

To: ALABAMA HIGHWAY DEPARTMENT



Report Date: 05/31/90

Attn: BUDDY COX

Lab number : 0105294 Sample number : HIWAY

Sample number: HIWAT
Sample matrix: WATER

COLLECTION INFORMATION

Date/Time/By: 05/25/90 11:25 COX Location : DANNELLY ANGB. 2-3-4

ADEM CENTRAL LABORATORY

- RESULTS REPORT - May 31, 1990

Lab#	Test	Result	UnitsDL*	Analdate
0105294	1,1,1,2-Tetrachloroetha 1,1,1-Trichloroethane 1,1,2,2-Tetrachloroetha	10.0000 469.7000 10.0000	ug/L	05/28/90 05/28/90 05/28/90
	1,1,2Trichloroethane 1,1-Dichloroethane	10.0000 10.0000	ug/L U ug/L U	05/28/90 05/28/90
-	1,1-Dichloroethylene 1,1-Dichloropropene	10.0000 10.0000 10.0000	ug/L U	05/28/90 05/28/90 05/28/90
	1,2,3-Trichlorobenzene 1,2,3-Trichloropropane 1,2,4-Trichlorobenzene	10.0000		05/28/90 05/28/90
	1,2,4-Trimethylbenzene 1,2-Dicholoethane	10.0000	ug/L U	05/28/90 05/28/90
	1,2-Dichloropropane 1,3,5-Trimethylbenzene 1,3-Dichloropropane	10.0000 10.0000 10.0000	ug/L U	05/28/90 05/28/90 05/28/90
	1,3-Dichloropropene 2,2-Dichloropropane	10.0000	ug/L U	05/28/90 05/28/90
	Tetrachloroethylene Bromobenzene	10.0000	ug/L U	05/28/90 05/28/90
•	Bromochloromethane Bromodichloromethane Benzene	10.0000 10.0000 - 10.0000	— ·	05/28/90 05/28/90 05/28/90

* U denotes results less than the instrument detection limit.

LABORATORY:	/ Montgomery	() Mobile	() Birmingham
Sample Type:	Potable Water [] Landfill L Surface Water [] Hazardous Soil/Sediment [] Groundwate Wastewater [] Waste (Speci	Wastesite [] Ignitab	vity [] Container P
Source Da	well, ANGB US80		
Location2-3-	4 BTEX - Water		
() Discharge	from	to	
Comments Skon	(Point Source) Kerosur obor- Kerosur of hones	Preservative(s)	(Receiving Water)
pH	D.O Sp. Cond	Salinity T	urb
Date	Value Date Value		
(mg/1)	(mg/1)	Date Valu (mg/l)	e Date Value (mg/l)
Acid	Pheno1		Mn
ALK	POP	Ag	Na
BOD ₅ -	(s ⁼)	As	Ni ·
(C1 ⁻)	(so ₄ =)	Ba	Pb
COD	TSS	Ca	Pt.
CN¯	TDS	Cd	Sb
(F ⁻)	TFS	Cr ^T	Se
Hard	TKN	Cr ⁺⁶	Zn
NH3-N	тос	Cu	Other
NO ₃ -N _	TON	Fe	BIEX-Vol
NO ₂ -N	TS	Нд	
0 & G	vss	Mg	
			F. Coli
B.C.	D. S.C. 11:25 1/2	16. RS/	7 (/ 24 /2/
SAMPLE COLLECT			(Signature) DATE/TIME
	•		
RECEIVED BY	(Signature) DATE/TI	ME RELINQUISHED BY	(Signature) DATE/TIME
RECEIVED BY,	(Signature) DATE/TIME Signature) DATE/TIME AB BY (Signature) DATE/TIME AB CONTROL OF THE SIGNATURE DATE/TIME AB CONTROL OF THE SIGN)	(Signature) DATE/TIME 0/05294
SEND REPORT TO			J •

TSS - Total Non-Filtrable Residue

TDS = Total Filtrable Residue

ADEM Form 68 5/83

TS = Total Residue

- SAMPLE ANALYSIS REPORT - 06/22/90

To: ALABAMA HIGHWAY DEPARTMENT

RECEIVED
Addra
Fict office
Monte.

Report Date: 06/22/90

Attn: BUDDY COX

Lab number : 0105305 Sample number : HIWAY

Sample matrix : SOIL

COLLECTION INFORMATION

Date/Time/By: 05/25/90 11:25 COX Location : DANNELLY ANGB. 3-1-3/5

ADEM CENTRAL LABORATORY

- RESULTS REPORT - June 22, 1990

Lab#	Test	Result	UnitsDL*	Analdate
0105305	1.2,4,-Trichlorobenzene	0.33	ug/L U	05/28/90
	1,2-Dichlorobenzene		ug/L U	05/28/90
	1,2-Diphenylhydrazine		ug/L U	05/28/90
	1.3-Dichlorobenzene		ug/L U	05/28/90
	1.4-Dichlorobenzene		ug/L U	05/28/90
	2.3.7.8-Tetrachlorodibe		ug/L U	05/28/90
	2.4-Dinitrotoluene		ug/L U	05/28/90
	2.6-Dinitrotoluene	0.33	ug/L U	05/28/90
	2-Chloronaphthalene	0.33	ug/L U	05/28/90
	3.3'-Dichlorobenzidine	0.33	ug/L U	05/28/90
	4-Bromophenyl phenyl eh	0.33	ug/L U	05/28/90
	4-Chlorophenyl phenyl e	0.33	ug/L U	05/28/90
	Acenaphthalene	0.33	ug/L U	05/28/90
	Acenaphthene	0.33	ug/L U	05/28/90
	Silver-EP	0.05	mg/L U	06/13/90
	Anthroene	0.33	ug/L U	05/28/90
•	Arsenic-EP	0.01	mg/L U	06/13/90
	Benzo(a)anthacene	0.33	ug/L U	05/28/90
	Barium-EP	0.50	mg/L U	06/13/90
	Benzo(a)pyrene	0.33	ug/L U	05/28/90
	Benzo(b)fluoanthene	0.33	ug/L U	05/28/90
	Butyl benzyl phthalate	0.33	ug/L U	05/28/90

U denotes results less than the instrument detection limit.

Lab#	Test	Result	UnitsDL*	Analdate
0105305	Bis (2-chlororthyl) eth	0.33	ug/L U	05/28/90
	Bis(2-chloroethoxy)meth		ug/L U	05/28/90
	Bis (2-Chloroisopropyl)		ug/L U	05/28/90
	Bis(2-ethylhexyl)phthal		ug/L	05/28/90
	Benzo(g,h,i)pervlene		ug/L U	05/28/90
	Benzidine		ug/L U	05/28/90
	Benzo(k)fluoranthene		ug/L U	05/28/90
	Cadmium-EP		mg/L U	06/13/90
	Chromium-EP		mg/L U	06/13/90
	Chrysene		ug/L U	05/28/90
	Dibenzo(a,h)anthracene		ug/L U	05/28/90
	Dibutyl phthalate		ug/L U	05/28/90
	Diethyl phthalate		ug/L U	05/28/90
	Dimethylphthalate		ug/L U	05/28/90
	Di-n-octyl phthalate		ug/L U	05/28/90
	Fluoranthene		ug/L U	05/28/90
•	Fluorene		ug/L U	05/28/90
	Hexachlorobutadiene		ug/L U	05/28/90
	Hexachlorobenzene		ug/L U	05/28/90
	Hexachlorocyclopentadie	0.33	IJ	05/28/90
	Hexachloroethane	0.33	ug/L U	05/28/90
	Mercury-EP		mg/L U	06/13/90
	Isophrene		ug/L U	05/28/90
	Indeno(1,2,3-cd)pyrene		ug/L U	05/28/90
	Naphthalene	0.33	ug/L U	05/28/90
	Nitrobenzene	0.33	ug/L U	05/28/90
	N-nitroso-di-n-propylam		ug/L U	05/28/90
	N-nitrosodimethylamine	0.33	ug/L U	05/28/90
	N-nitrosodiphenylamine	9.33	ug/L U	05/28/90
	Pyrene		ug/L U	05/28/90
	Phenanthrene		ug/L U	05/28/90
	Lead-EP		mg/L U	06/13/90
	Lead in Soil	61.7	ug/g	06/13/90
	Selenium-EP	0.01	mg/L U	06/13/90

^{*} U denotes results less than the instrument detection limit.

LABORATORY:	✓ Montgomery	() Mobile	() Birmingham
Surface	Water [] Landfill Leach Water [] Hazardous Wast diment [] Groundwater ter [] Waste (Special H	tesite [ˈ] Ignitabilii [] Corrosivity	Container P
Source <u>3-/-3/5</u>	Danvelly ANGB-	US80)	
Location 3-1-3/5			
() Discharge from		to .	
	(Point Source)		(Receiving Water)
Comments		Preservative(s)	
D.O	Sp. CondPARA	Salinity Turb	
Date Value (mg/l)	Date Value (mg/l)	Date Value	Date Value
Acid	Phenol	(mg/l) Al	(mg/l) Mn
ALK	PO, -P	Ag	Na
BOD ₅	(S ⁼)	As	Ni Ni
(C1 ⁻)	(so ₄ =)	Ba	Pb
COD	TSS	Ca	Pt
CN	TDS	Cd	Sb
(F ⁻)	TFS	Cr ^T	Se
llard	TKN	Cr ⁺⁶	Zn
NH ₃ -N	TOC	Cu	Other
NO ₃ -N	TON	Fe	Base Newfiels
NO ₂ -N	TS	Нд	Pb (Tof + EP)
O & G	VSS	Mg	
	5 . //		F. Coli.
SAMPLE POLLECTED BY	B. <u>5</u> (1/.25 5/25/90 (Signature) DATE/TIME	RÉLINQUISHED BY (S	Ignature) DATE/TIME
as yourster bi	(organie) Dait/IIME	verindatouen bi (2)	nature) DATE/TIME
RECEIVED BY	(Signature) DATE/TIME	RELINQUISHED BY (S	ignature) DATE/TIME
RECEIVED BY	(Signature) DATE/TIME	-	ignature) DATE/TIME 5305
RECEIVED IN LAB BY	(Signature) DATE/TIME	LABORATORY I.D. NO.	<u> </u>
SEND REPORT TO:	ox (Mat + Test)	· · · · · · · · · · · · · · · · · · ·	

- SAMPLE ANALYSIS REPORT -06/13/90

To: ALABAMA HIGHWAY DEPARTMENT

Nonig.

Attn: BUDDY COX

Lab number : 0105301 Sample number : HIWAY . Sample matrix : SOIL

Report Date: 06/13/90

COLLECTION INFORMATION Date/Time/By: 05/25/90 11:20 COX Location : DANNELLY ANGE, 3-2-3/5

> ADEM CENTRAL LABORATORY - RESULTS REPORT -

June 13, 1990

Lab#	Test	Result	UnitsDL*	Analdate
0105301	1,2,4Trichlorobenzene		ug/L U ug/L U	05/28/90
	1,2-Diphenylhydrazine		ug/L U	05/28/90 05/28/90
	1.3-Dichlorobenzene		ug/L U	05/28/90
	1,4-Dichlorobenzene		ug/L U	05/28/90
	2,3,7,8-Tetrachlorodibe		ug/L U	05/28/90
	2.4-Dinitrotoluene		ug/L U	05/28/90
	2.6-Dinitrotoluene		ug/L U	05/28/90
	2-Chloronaphthalene		ug/L U	05/28/90
	3.3'-Dichlorobenzidine		ug/L U	05/28/90
	4-Bromophenyl phenyl eh		ug/L U	05/28/90
	4-Chlorophenyl phenyl e		ug/L U	05/28/90
	Acenaphthalene	0.33	ug/L U	05/28/90
•	Acenaphthene	0.33	ug/L U	05/28/90
	Silver-EP	0.05	mg/L U	06/11/90
	Anthroene	0.33	ug/L U	05/28/90
•	Arsenic-EP	0.01	mg/L U	06/11/90
	Benzo(a)anthacene		ug/L U	05/28/90
	Barium-EP		mg/L U	06/11/90
	Benzo(a)pyrene		ug/L U	05/28/90
	Benzo(b)fluoanthene		ug/L U	05/28/90
	Butyl benzyl phthalate	0.33	ug/L U	05/28/90

^{*} U denotes results less than the instrument detection limit.

Lab#	Test	Result	UnitsDL*	Analdate
0105301	Bis (2-chlororthyl) eth	0.33	ug/L U	05/28/90
	Bis(2-chloroethoxy)meth		ug/L U	05/28/90
	Bis (2-Chloroisopropyl)	0.33	ug/L U	05/28/90
	Bis(2-ethylhexyl)phthal	0.33	ug/L U	05/28/90
	Benzo(g.h,i)perylene	0.33	ug/L U	05/28/90
	Benzidine	0.33	ug/L U	05/28/90
	Benzo(k)fluoranthene	0.33	ug/L U	05/28/90
	Cadmium-EP	0.01	mg/L U	06/11/90
	Chromium-EP	0.05	mg/L U	06/11/90
	Chrysene	0.33	ug/L U	05/28/90
	Dibenzo(a,h)anthracene	0.33	ug/L U	05/28/90
	Dibutyl phthalate	0.33	ug/L U	05/28/90
	Diethyl phthalate	0.33	ug/L U	05/28/90
	Dimethylphthalate	0.33	ug/L U	05/28/90
	Di-n-octyl phthalate		ug/L U	05/28/90
	Fluoranthene	0.33	ug/L U	05/28/90
	Fluorene	0.33	ug/L U	05/28/90
	Hexachlorobutadiene		ug/L U	05/28/90
	Hexach Lorobenzene		ug/L/U	05/28/90
	Hexachlorocyclopentadie	0.33	Ū	05/28/90
	Hexachloroethane		ug/L U	05/28/90
	Mercury-EP		mg/L U	06/11/90
	Isophrone		ug/L U	05/28/90
	Indeno(1,2,3-cd)pyrene		ug/L U	05/28/90
	Naphthalene		ug/L U	05/28/90
•	Nitrobenzene		ug/L U	05/28/90
	N-nitroso-di-n-propylam		ug/L U	05/28/90
	N-nitrosodimethylamine		ug/L U	05/28/90
	N-nitrosodiphenylamine		ug/L U	05/28/90
	Pyrene		ug/L U	05/28/90
	Phenanthrene		ug/L U	05/28/90
	Lead-EP		mg/L U	06/11/90
	Lead in Soil		ug/g	06/11/90
	Selenium-EP	0.01	mg/L U	06/11/90

^{*} U denotes results less than the instrument detection limit.

LABORATORY:	(Montgomery	()	Mobile	() Birmingham	1
Sample Type:	Surface Water [] Haza Soil/Sediment [] Grou	fill Leachate rdous Wastesite ndwater e (Special Handling)	[] Ignitabil [] Corrosivi	ty [] Container	ĺ
Source	annelly ANGB W	1880			
Location 3	2-3/5				
() Discharge			to		
Comments	(Point S		ervative(s)	(Receiving Water)	,
pH	D.O Sp. Cond.	·			
Date (mg/l)	Value Date (mg/1)	Value Date	value (mg/l)	Date Valu	16 4
Acid	-		Al	(mg/1) Mn	
ALK	PO, -P		Ag	· Na	
BOD ₅	(s ⁼)		As	Ni Ni	_ [
(C1 ⁻)	(SO ₄ =)		Ba	Pb	
COD	TSS		Ca	Pt Pt	_ 1
CN	TDS		Cd		
(F ⁻)	TFS		Cr ^T	Se	
Hard	TKN			Zn	
NH_3-N	TOC		Cu	Other	
NO_3-N	TON		— Fe	Bax Newfood	
NO ₂ -N	TS		Hg	Rh (Tot JEP7	フ
0 & G	vss		 Mg		
				F. Coli	
B. Cox	7516/ 51	//:20 //s=/sa		A.Gx 2:00 5/2	(=)
SAMPLE COLLEC	TED BY (Signature) D	ATE/TIME RELIN	QUISHED BY (Signature) DATE/TIME	3 <i>/7</i>
RECEIVED BY	(Signature) D	ATE/TIME RELIN	QUISHED BY (Signature) DATE/TIME	2
RECEIVED BY	(Signature) D	ATE/TIME RELIN	QUISHED BY (Signature) DATE/TIME	2
Delots		2:00	0/	05301	
RECEIVED IN L	- " ~		RATORY I.D. NO.		
SEND REPORT T	: Cox (Mat. +)	Test / Hijhung	<i></i>		

ADEM CENTRAL LABORATORY

- SAMPLE ANALYSIS REPORT -05/31/90

To: ALABAMA HIGHWAY DEPARTMENT

Jiii 1990

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RECEIVED

Adom Field Cifice Montg

Attn: BUDDY COX

Lab number : 0105293

Report Date: 05/31/90

Sample number : HIWAY Sample matrix : WATER

COLLECTION INFORMATION

Date/Time/By: 05/25/90 11:50 COX Location : DANNELLY ANGB, 3-4-4

> ADEM CENTRAL LABORATORY - RESULTS REPORT -

May 31, 1990

0105293 1.1.1.2-Tetrachloroetha 10.0000 ug/L U 05/28 1.1,1-Trichloroethane 10.0000 ug/L U 05/28 1,1,2.2-Tetrachloroetha 10.0000 ug/L U 05/28 1,1,2Trichloroethane 10.0000 ug/L U 05/28	date
1.1-Dichloroethane 10.0000 ug/L U 05/20 1.1-Dichloropropene 10.0000 ug/L U 05/20 1.2.3-Trichlorobenzene 10.0000 ug/L U 05/20 1.2.3-Trichloropropane 10.0000 ug/L U 05/20 1.2.4-Trichlorobenzene 10.0000 ug/L U 05/20 1.2.4-Trimethylbenzene 10.0000 ug/L U 05/20 1.2-Dichloropropane 10.0000 ug/L U 05/20 1.2-Dichloropropane 10.0000 ug/L U 05/20 1.3.5-Trimethylbenzene 10.0000 ug/L U 05/20	8/90 8/90 8/90 8/90 8/90 8/90 8/90 8/90
1,3-Dichloropropane 10.0000 ug/L U 05/20 1,3-Dichloropropene 10.0000 ug/L U 05/20 2,2-Dichloropropane 10.0000 ug/L U 05/20 Tetrachloroethylene 10.0000 ug/L U 05/20	8/90 8/90
Bromobenzene 10.0000 ug/L U 05/2 Bromochloromethane 10.0000 ug/L U 05/2 Bromodichloromethane 10.0000 ug/L U 05/2	:8/90 :8/90

^{*} U denotes results less than the instrument detection limit.

Lab#	Test	Result	UnitsDL*	Analdate
0105293	Bromomethane	10.0000		05/28/90
	cis-1,2-Dichloroethylen	10.0000		05/28/90
	Chlorobenzene	10.0000		05/28/90
	Chlorodibromomethane	10.0000		05/28/90
	Chloroethane	10.0000		05/28/90
	Bromoform	10.0000		05/28/90
	Chloroform	10.0000	ug/L U	05/28/90
	Chloromethane	10.0000	ug/L U	05/28/90
	Carbon Tetrachloride	10.0000	ug/L U	05/28/90
	Dibromomethane	10.0000	ug/L U	05/28/90
	Dichlorofluoromethane	10.0000	ug/L U	05/28/90
	Dichloromethane	10.0000	ug/L U	05/28/90
	Ethylbenzene	10.0000	ug/L U	05/28/90
	Fluorotrichloromethane	10.0000	ug/L U	05/28/90
	Hexachlorobenzene	10.0000		05/28/90
	Isopropylbenzene	10.0000		05/28/90
	m-Dichlorobenzene	10.0000		05/28/90
	m-Xylene	10.0000		05/28/90
	Naphthalene	10.0000		05/28/90
	n-Butylbenzene	10.0000		05/28/90
	n-Propylbenzene	10.0000		05/28/90
	o-Chlorotoluene	10.0000		05/28/90
	o-Dichlorobenzene	10.0000	ug/L U	05/28/90
	o-Xylene	10.0000	ug/L U	05/28/90
	p-Chlorotoluene	10.0000	ug/L U	05/28/90
	p-Dichlorobenzene	10.0000	ug/L U	05/28/90
	p-Isopropyltoluene	10.0000	ug/L U	05/28/90
	p-Xylene	10.0000	ug/L U	05/28/90
	Secbutylbenzene	10.0000	ug/L U	05/28/90
	Styrene	10.0000	ug/L U	05/28/90
	t-1,2Dichloroethane	10.0000	ug/L U	05/28/90
	Tertbutylbenzene	10.0000	ug/L U	05/28/90
	Trichloroethylene	10.0000	ug/L U	05/28/90
	Toluene	10.0000		05/28/90
	Vinyl Chloride	10.0000	ug/L U	05/28/90

^{*} U denotes results less than the instrument detection limit.

LABORATORY:	<i>∞</i> ма	ontgomery	() Mobile	() Birmingham
Sample Type:	Surface Water Soil/Sediment	Hazardous Waste Groundwater	esite [] Ignitabi	ity [] Container P
Source Da	well ANG	В		
Location 3-	4-4 Weste			
() Discharge	from	(Point Source)	to	
Comments			_ Preservative(s)	(Receiving Water)
]	D.U :	PARAM	SalinityTu ETERS	rb
Date (mg/1)	Value Dat	e Value (mg/1)	Date Value (mg/l)	Date Value (mg/l)
Acid_		Phenol	A1	Mn
ALK _		PO ₄ -P	Ag	Na
BOD ₅ -		(s ⁼)	As	Ni
(C1 ⁻)		(SO ₄ =)	Ba	Pb
COD _		TSS	Ca	Pt
CN		TDS	Cd	Sb
(F ⁻)		TFS	Cr ^T	Se
Hard _		TKN	Cr ⁺⁶	Zn
NH ₃ -N -		TOC	Cu	Other
NO ₃ -N -		TON	Fe	BTEX - vol
NO ₂ -N _		TS	Нд	-
0 & G		vss	Mg	
B. S. U	B. 5.4	 11:505/15 to	4	F. Coli
SAMPLE COLLECT	ED BY (Signat	ure) DATE/TIME		2:05 5/25/90 (Signature) DATE/TIME
RECEIVED BY	(Signat	ure) DATE/TIME	RELINQUISHED BY	(Signature) DATE/TIME
RECEIVED BY	(Signat	ure) DATE/TIME	RELINQUISHED BY	(Signature) DATE/TIME
Melin	Odener			
RECEIVED IN LA			LABORATORY I.D. NO	5293
	. 8	ex Materials +		
	7	MANIE O	1-01	

TSS = Total Non-Filtrable Residue

ADEM CENTRAL LABORATORY

- SAMPLE ANALYSIS REPORT - 05/31/90

To: ALABAMA HIGHWAY DEPARTMENT



Report Date: 05/31/9

Attn: BUDDY COX

Lab number : 0105295

Sample number : HIWAY Sample matrix : SOIL

COLLECTION INFORMATION

Date/Time/By: 05/25/90 11:50 COX Location : DANNELLY ANGB, 3-3-2

ADEM CENTRAL LABORATORY
- RESULTS REPORT -

May 31, 1990

Lab#	Test	Result	UnitsDL*	Analdate
0105295	1,1,1,2-Tetrachloroetha	0.0500	ug/g U	05/29/90
	1,1,1-Trichloroethane	0.0500		05/29/90
	1,1,2,2-Tetrachloroetha	0.0500		05/29/90
	1,1,2Trichloroethane	0.0500		05/29/90
	1,1-Dichloroethane	0.0500	ug/g U	05/29/90
	1,1-Dichloroethylene	0.0500	ug/g U	05/29/90
	1,1-Dichloropropene	0.0500	ug/g U	05/29/90
	1,2,3-Trichlorobenzene	0.0500	ug/g U	05/29/90
	1,2,3-Trichloropropane	0.0500	ug/g U	05/29/90
	1.2.4-Trichlorobenzene	0.0500	ug/g U	05/29/90
	1,2,4-Trimethylbenzene	0.0500	ug/g U	05/29/90
	1,2-Dicholoethane	0.0500	ug/g U	05/29/90
	1,2-Dichloropropane	0.0500	ug/g U	05/29/90
	1,3,5-Trimethylbenzene	0.0500	ug/g U	05/29/90
	1,3-Dichloropropane	0.0500	ug/g []	05/29/90
•	1,3-Dichloropropene	0.0500	ug/g U	05/29/90
	2.2-Dichloropropane	0.0500	ug/g U	05/29/90
	Tetrachloroethylene	0.0500	ug/g U	05/29/90
	Bromobenzene	0.0500	ug/g U	05/29/90
	Bromochloromethane	0.0500	ug/g U	05/29/90
	Bromodichloromethane	0.0500	ug/g U	05/29/90
	Benzene	0.0500	ug/g U	05/29/90

U denotes results less than the instrument detection limit.

Lab#	Test	Result	UnitsDL:*	Analdate
Lab# 	Bromomethane cis-1,2-Dichloroethylen Chlorobenzene Chlorodibromomethane Chloroethane Bromoform Chloroform Chloromethane Carbon Tetrachloride Dibromomethane Dichlorofluoromethane Dichlorofluoromethane Ethylbenzene Fluorotrichloromethane Hexachlorobenzene Isopropylbenzene m-Dichlorobenzene m-Dichlorobenzene m-Xylene Naphthalene n-Butylbenzene	0.0500 0.0500 0.0500 0.0500 0.0500 0.0500 0.0500 0.0500 0.0500 0.0500 0.0500 0.0500 0.0500 0.0500	UnitsDL* ug/g U	Analdate 05/29/90
	n-Butylbenzene n-Propylbenzene o-Chlorotoluene o-Dichlorobenzene o-Xylene p-Chlorotoluene p-Dichlorobenzene p-Isopropyltoluene p-Xylene Secbutylbenzene Styrene t-1,2Dichloroethane Tertbutylbenzene Trichloroethylene Toluene Vinyl Chloride	0:0500 0.0500 0.0500 0.0500 0.0500 0.0500 0.0500 0.0500 0.0500 0.0500 0.0500	ug/g U	05/29/90 05/29/90 05/29/90 05/29/90 05/29/90 05/29/90 05/29/90 05/29/90 05/29/90 05/29/90 05/29/90 05/29/90 05/29/90 05/29/90

^{*} U denotes results less than the instrument detection limit.

Ai-Way

LABORATORY:	Montg	omery	() Mobile		() Birmi	ngham
Sample Type:	Surface Water []	Groundwater	stesite [] I [] C	gnitabili Orrosivit	action [] ty [] y []	Grab
Source	anuelly ANGB	W80				
Location 3-	3-2					
() Discharge			t	.0		
		int Source)			(Receiving	
Comments						
pH	D.O Sp.	Cond. PAR	Salinity RAMETERS	Turb		
Date		Value	Date	Value	Date	Value
(mg/1)		mg/1)	(mg/1)	. (mg/1)
Acid	•	henol	A1			Mn
ALK		0 ₄ -P	Ag			Na
BOD ₅ -		(S ⁼)	As			Ni
(C1 ⁻)		so ₄ =)	Ba			Pb
COD .		TSS	Ca			Pt
CN		TDS	Cd			Sb
		TFS	Cr [†]			Se
Hard Hard		TKN	Cr ⁺⁶			Zn
NH ₃ -N		TOC	Cu		Other	
NON		TON	Fe		BIE	Vol
NO ₂ -N _		TS	Нд			
		vss	Mg			<u></u>
PIO 6	5//	•			F. Coli.	
SAMPLE COLLECT	ED BY (Signature	// <i>SOB/25/9</i> 0) DATE/TIME	B. GC RELINQUISHE	D BY (S	ignature)	2:00 5/25/9 DATE/TIME
RECEIVED BY	(Signature) DATE/TIME	RELINQUISHE	D BY (S	ignature)	DATE/TIME
RECEIVED BY RECEIVED IN LA	(Signature	2:00	RELINQUISHE		ignature) 01052	DATE/TIME
SEND REPORT TO	\sim $^{\circ}$	DATE/TIME	LABORATORY	I.D. NO.		
CENT RELORI IC	· - or / Ma	1 + 1 - 31			*	

- SAMPLE ANALYSIS REPORT -06/22/90

To: ALABAMA HIGHWAY DEPARTMENT

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JU1 - 1

LECENED Adom Note Office Monte.

Attn: BUDDY COX

Lab number : 0105304

Report Date: 06/22/90

Sample number : HIWAY Sample matrix : SOIL

COLLECTION INFORMATION ...

Date/Time/By: 05/25/90 12:15 COX Location : DANNELLY ANGB. 4-1-3/5

ADEM CENTRAL LABORATORY

- RESULTS REPORT

June 22, 1990

Lab#	Test	Result	UnitsDL+	Analdate
Lab# 0105304	1.2.4Trichlorobenzene 1.2-Dichlorobenzene 1.2-Diphenylhydrazine 1.3-Dichlorobenzene 1.4-Dichlorobenzene 2.3.7.8-Tetrachlorodibe 2.4-Dinitrotoluene 2.6-Dinitrotoluene 2-Chloronaphthalene 3.3'-Dichlorobenzidine 4-Bromophenyl phenyl eh 4-Chlorophenyl phenyl eh 4-Chlorophenyl phenyl e Acenaphthalene Acenaphthalene Acenaphthene Silver-EP Anthroene Arsenic-EP Benzo(a)anthacene Barium-EP	0.33 0.33 0.33 0.33 0.33 0.33 0.33 0.33	UnitsDL* ug/L U	Analdate 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 06/13/90 05/28/90 06/13/90 06/13/90
	Benzo(a)pyrene Benzo(b)fluoanthene Butyl benzyl phthalate	0.33 0.33	ug/L U ug/L U ug/L U	05/28/90 05/28/90 05/28/90

^{*} U denotes results less than the instrument detection limit.

Lab# 	Test	Result	UnitsDL*	Analdate
Lab# 0105304	Bis (2-chlororthyl) eth Bis (2-chloroethoxy)meth Bis (2-chloroisopropyl) Bis(2-ethylhexyl)phthal Benzo(g,h,i)perylene Benzidine Benzo(k)fluoranthene Cadmium-EP Chromium-EP Chromium-EP Chrysene Dibenzo(a,h)anthracene Dibutyl phthalate Diethyl phthalate Dimethylphthalate Di-n-octyl phthalate Fluoranthene Fluorene Hexachlorobutadiene Hexachlorobenzene Hexachloroethane Mercury-EP Isophrone Indeno(1,2,3-cd)pyrene Naphthalene Nitrobenzene N nitropo-di n propylam N-nitrosodimethylamine N-nitrosodimethylamine Pyrene	0.33 0.33 0.33 0.33 0.33 0.33 0.33 0.33	Unitable ug/L U ug/L	Analdate 05/28/90
	Phenanthrene Lead-EP Lead in Soil Selenium-EP	0.33 0.20 23.7	ug/L U mg/L U	05/28/90 05/28/90 06/13/90 06/13/90 06/13/90

^{*} U denotes results less than the instrument detection limit.

Hiway

LABORATORY:	Montg	omery	() Mobile	() Birmingham
Sample Type:	Potable Water [] Surface Water [] Soil/Sediment [] Wastewater []	Landfill Leach Hazardous Wass Groundwater Waste (Special H	esite [] Ignita [] Corros	Extraction Composite bility Grab Container P Coulty G
Source Da	unelly ANGB-	Fence Lin	e	
Location 4	I			
() Discharge	from	int Source)	to	
				(Receiving Water)
Comments	· · · · · · · · · · · · · · · · · · ·		_ Preservative(s)	
рН	D.OSp.		Salinity METERS	Turb.
Date (mg/l)	Value Date	Value mg/l)	Date Val	ue Date Value (mg/l)
Acid		henol	Al	Mn
ALK		O, -P	Ag	Na
BOD		(S ⁼)	As	Ni
(C1 ⁻)		so ₄ =)	Ba	(Pb)
COD		rss	Ca	Pt
CN		rds	Cd	Sb
(F ⁻)		rfs	Cr ^T	
Hard _		rkn	Cr ⁺⁶	Zn
NH3-N _		гос	Cu	Other
NO ₃ -N	· · · · · · · · · · · · · · · · · · ·		Fe	Base Nuctrals
NO ₂ -N _		TS	Hg	Lead (Tot 4EP)
0 & G _		vss	Mg	
l BEA	- 344//	a		F. Coli.
SAMPLE COLLECT	ED BY (Signature	5/25/90 /2/5 DATE/TIME	RELINQUISHED BY	(Signature) DATE/TIME
			REDINGOISHED BI	(Sygnacure) DATE/TIME
RECEIVED BY	(Signature) DATE/TIME	RELINQUISHED BY	(Signature) DATE/TIME
RECEIVED BY	(Şignature) DATE/TIME	RELINQUISHED BY	(Signature) DATE/TIME
W. M.		7:00		0105304
RECEIVED IN LA	B BY (Signature		LABORATORY I.D.	NO.
SEND REPORT TO	: Cox	(Mat & Tes)	() 242-6257	

- SAMPLE ANALYSIS REPORT - 06/22/90

To: ALABAMA HIGHWAY DEPARTMENT



Attn: BUDDY COX

Lab number : 0105306 Sample number : HIWAY Sample matrix : SOIL

Report Date: 06/22/9/

COLLECTION INFORMATION

Date/Time/By: 05/25/90 12:40 COX Location : DANNELLY ANGB, 4-3-3/5

ADEM CENTRAL LABORATORY
- RESULTS REPORT -

June 22, 1990

Lab#	Test	Result	UnitsDL*	Analdate
0105306	1,2,4,-Trichlorobenzene 1,2-Dichlorobenzene 1,2-Diphenylhydrazine 1,3-Dichlorobenzene 1,4-Dichlorobenzene 2,3,7,8-Tetrachlorodibe 2,4-Dinitrotoluene 2,6-Dinitrotoluene 2-Chloronaphthalene 3,3'-Dichlorobenzidine 4-Bromophenyl phenyl el A-Chlorophenyl phenyl el Acenaphthalene Acenaphthene Silver-EP Anthroene Arsenic-EP Benzo(a)anthacene Barium-EP Benzo(b)fluoanthene Butyl benzyl phthalate	0.33 0.33 0.33 0.33 0.33 0.33 0.33 0.33	ug/L U	05/28/20 05/28/20 05/28/90 05/28/90

^{*} U denotes results less than the instrument detection limit.

Lab#	Test	Result	UnitsDL*	Analdate
Lab# 0105306		0.33 0.33 0.33 0.33 0.33 0.33 0.33 0.33	ug/LU	05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 05/28/90 06/13/90 06/13/90 05/28/90
	Lead in Soil Selenium-EP	25.1	mg/L U ug/d mg/∐ U	06/13/90 06/13/90 06/13/90

^{*} U denotes results less than the instrument detection limit.

LABORATORY:	Montgomery	() Mobile () Birmingham	1
Sample Type:	Surface Water [] Hazardous Was Soil/Sediment Groundwater	chate [] Toxic Extraction M. Composite	. 7
Source Dans	celly ANGO 4580		1
Location 4			_
() Discharge	from	to.	
	(Point Source)	(Receiving Water)	_
Comments			
pH	D.O Sp. CondPAR/	_ Salinity Turb	
	Value Date Value	Data v.1 -	
(mg/1)	(mg/1)	Date Value Date Value (mg/l) (mg/l)	•
Acid _	Phenol	Al Mn	1
ALK _	PO ₄ -P	Ag Na	
BOD ₅ -	(s ⁼)	As Ni	ı
(C1 ⁻)_	(so ₄ =)	Ba	
COD _	TSS	Ca Pt	•
CN	TDS	Cd Sb	
(F¯)_	TFS	Cr ^T Se	_
Hard _	TKN	Cr ⁺⁶ Zn	
NH ₃ -N _	TOC	Cu Other	
NO ₃ -N _	TON	Fe Base New /Pb	1
NO ₂ -N _	TS	HgTotals	, =
O & G	vss	Mg	1
B5///		F. Coli	
SAMPLE COLLECTI	ED BY (Signature) DATE/TIME	RELINQUISHED BY (Signature) DATE/TIME	
·	. /	(Orginature) DATE/TIME	
RECEIVED BY	(Signature) DATE/TIME	RELINQUISHED BY (Signature) DATE/TIME	-1
RECEIVED BY ARECEIVED IN LAI	(Signature) DATE/TIME Shure 5/25 2:00 By (Signature) DATE/TIME	RELINQUISHED BY (Signature) DATE/TIME 005306 LABORATORY I.D. NO.	
SEND REPORT TO:	Cox (Map. o Test)	<u>.</u>	
	242		

TSS - Total Non-Filtrable Residue

TOS = Total Filtrable Residue

ADEM Form 68 5/83

TS = Total Residue

Sample matrix : SOIL

COLLECTION INFORMATION

Date/Time/By: 05/25/90 12:35 COX Location : DANNELLY ANGB, 4-2-2

ADEM CENTRAL LABORATORY
- RESULTS REPORT -

May 31, 1990

Lab#	Test	Result	UnitsDL*	Analdate
0105296	1,1,1,2-Tetrachloroetha 1,1,1-Trichloroethane 1,1,2,2-Tetrachloroetha 1,1,2Trichloroethane 1,1-Dichloroethane 1,1-Dichloroethylene 1,1-Dichloropropene 1,2,3-Trichlorobenzene 1,2,3-Trichloropropane 1,2,4-Trimethylbenzene 1,2-Dichloropropane 1,2-Dichloropropane 1,3-Trimethylbenzene 1,3-Dichloropropane	0.0500 0.0500 0.0500 0.0500 0.0500 0.0500 0.0500 0.0500 0.0500 0.0500 0.0500 0.0500 0.0500 0.0500 0.0500 0.0500	ug/g U ug/g U ug/g U ug/g U ug/g U	05/29/90 05/29/90

* U denotes results less than the instrument detection limit.

— ·— — — — — — — — — — — — — — — — — —		
Toluene	0.0500 ug/g U 05/29/90	
Vinyl Chloride	0.0500 ug/g U 05/29/90	

^{*} U denotes results less than the instrument detection limit.

LABORATORY:	Montgomery	() Mobile	() Birmingham
Sample Type:	i) Deacite	esite [] Ignitab [] Corrosiv	vity [] Container P i
Source Jan	nely ANGB US80		/ =
Location 4	× ×		
() Discharge	from (Point Source)	to	(Receiving Water)
Comments		Preservative(s)	(Receiving water)
рН	D.OSp. CondS	Salinity To	
	Value Date Value	Date Value	Date Value
(mg/1)	(mg/1)	(mg/l)	(mg/1)
Acid _	Phenol	A1	Mn
BOD ₅ -	PO ₄ -P	Ag	Na
(C1 ⁻)		As	Ni
COD	(SO ₄ =)	Ba	Pb
CN	TSS	Ca	Pt
(F ⁻)	TDS	Cd	Sb
Hard	TFS	Cr ^T	Se
	TKN	Cr ⁺⁶	Zn
NH ₃ -N _	TOC	Cu	Other
NON	TON	Fe	_ (3TEX-Vo/
NO ₂ -N _ O & G	TS	Нд	
O & G _	vss	Mg	
7.4/	B. E. 4/ 12:35 5/25/90	B.4/	F. Coli
SAMPLÉ COLLECT	ED BY (Signature) DATE/TIME	RELINQUISHED BY	(Signature) DATE/TIME
RECEIVED BY	(Signature) DATE/TIME	RELINQUISHED BY	(Signature) DATE/TIME
RECEIVED BY	(Signature) DATE/TIME	010	(Signature) DATE/TIME
	O SHOULD DATE / TIME	LABORATORY 1.D. NO	
SEND REPORT TO	: _ Osx (Maf + Tes	<i>T</i>	

TSS - Total Non-Filtrable Residue
TDS - Total Filtrable Residue

TS = Total Residue
VSS = Volatile Residue

ALABAMA DEPARTMENT OF ENVIRONMENTAL MANAGEMENT



Guy Hunt Governor

Leigh Pegues, Director

1751 Cong. W. L.

June 19, 1990 **Dickinson Drive**

Montgomery, AL

36130

205/271-7700

TO:

Highway Department

Field Offices:

Unit 806, Building 8 225 Oxmoor Circle

Birmingham, AL 35209

205/942-6168

P.O. Box 953 Decatur, AL

35602

205/353-1713

2204 Perimeter Road Mobile, AL 36615 205/479-2336

MEMORANDUM

Buddy Cox

John Chitwood

SUBJECT:

FROM:

Laboratory Results

Attached are laboratory results from samples submitted to the ADEM

Central Laboratory for analysis by the Highway Department.

JC/mpt

Attachments

ALABAMA DEPARTMENT OF ENVIRONMENTAL MANAGEMENT



Guy Hunt Governor

Leigh Pegues, Director

1751 Cong. W. L. **Dickinson Drive** Montgomery, AL

36130

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35209

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P.O. Box 953 Decatur, AL 35602

205/353-1713

2204 Perimeter Road

Mobile, AL 36615

205/479-2336

MEMORANDUM

June 8, 1990

TO:

Buddy Cox

Highway Department

FROM:

John Chitwood JCC

SUBJECT:

Laboratory Results

Attached are laboratory results from samples submitted to the ADEM Central Laboratory for analysis by the Highway Department.

JC/mpt

Attachments

ALABAMA DEPARTMENT OF ENVIRONMENTAL MANAGEMENT



Guy Hunt Governor

Leigh Pegues, Director

1751 Cong. W. L. **Dickinson Drive** Montgomery, AL

36130

205/271-7700

ield Offices:

Mr. Buddy E. Cox

Montgomery, Alabama 36130

Unit 806, Building 8 225 Oxmoor Circle Birmingham, AL 35209

205/942-6168

P.O. Box 953 Decatur, AL 35602 205/353-1713

2204 Perimeter Road Mobile, AL 36615 205/479-2336

Materials & Tests Alabama Highway Department Fairgrounds Road

RE: Samples 0105293, 0105294, 0105295, 0105296, and 0105297

Dear Mr. Cox:

Attached please find results of samples submitted to the Department of Environmental Management Laboratory on May 5, 1990. Should you have any questions with regard to the results please contact Mr. Joe Marsh at 271-7980

Sincerely,

June 5, 1990

& John Williford

E. John Williford, Chief Field Operations Division

EJW/TM/km

Enclosure

cc: Mr. Joe Marsh

Tor

Atton

Our Lab # : 0103293 Your Sample ID: HIMAY Samule Matrix : WATER

Invoice No.: 0103191

Invoice Date: 05/30/

INVOICE SUMMARY: Test Charge: 150.00 Additional Chargest

150.0

ree

150.0·

TOTAL

ITEMIZED TEST CHARGES: Laby Test

0105293

h	F W 122	5. 65.14V
0105293	110CE	*** *** *******************************
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0105293	ores	
0103293	12005	
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0105293	CB	
0105293	11177E	େ ୦୦
0105293	MOCE,	
0105293	1112TE	
0105997	713 Str. 7	
0105593	CY	
0105293	59004	
0105093	TEECH	
0105293	E 6.	
0105253	SM	
0105293	YOL	
0103293	SEE	
0105293	1110P	
0105293	CE	

CHOLG

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			the management of the control of the
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	401/05/22	11	
	0105293	$\{af^{a}f^{c}\}$	
	03.05090	PERF	
	0105293	120106	
	0105293	This es	
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ľ	0105223	$_{ m CDG}$ $^{ m CG}$	•
	0400000	CHIRAL II	
	0105293	4. 油水 () ()	
	0105273	V.C	
	0105293	SIY	
	0105293	ECM	
	0105293	19905	
	0105293	DCFM	
	0105293	HER	
	0100293	IPE	•
	0105293	112176	
	0108293	17710	
	0105293	86	·
	0105293	ACLE	•
	0105293	11270E	
	0105293	C4 20	
	0105293	Fich	
	0105293	MEB	•

THATCHEE

Tor

Attr:

INVOICE SUMMARY:

Our Lab # . : 0165224 Your Sample ID: HIMAY

Sample Betrix : MATER

Test Charges 150.00

Invoice Date: 05/30/90

150.00

invoice Mo.: 0103192

Additional Charges:

\$ 150,00

ITEMIZED TEST CHARGES: Labii 110CE 0105294 0.00 1111CE 0100274 120CE 0105293 0X0105294 VC 0105294 CDEM 0105294 POCE 0105294 PX 0105294 gapor ' 0105294 $\mathbb{C}\mathbb{S}$ 0105274 FTCM 0105294 $\mathcal{C}(\mathbb{C}; \mathbb{I})$ 0105274 STY 0105293 13548 0105294 112TCE 0105254 $\Xi \mathbf{E}$ 0105294 Biri 0105294 TOL. 0105294 FEZ 0105294 CT 0105274 TIPEC 0105294 124700 0105294 f:F: 0100294 CHARG 0105294 122TF 0105294 136B 0105294 BOOM 0105294 150.00 TOE 0105274 0105294 CE CHOLG. 0105294

150.00

Lab#	Test	발표를 하고 선택하고 있는데 하면 보고 있는데 되었고 있다.
0105294		
		하다. 그리는 사람들이 하시아 그 말이라는 사이를 모든 어린을하다.
0105194	(† 15 †	그 이 집에 대한다. 그 나라는 바쁜 아이시에다. 글로그 모든
0105291	PCH 3	그리고 교통 하고 계계 이 그리고 있는 살은 것이 그를 이미리를 내워.
0905225		요즘 하게 많아들어 하셨다니까? 그렇게 되고 이번째 손님 되었다.
0:0579-	117211	입니다는 말하고 하고 말하게 되는 때 나는 것이 되어 있다.
0105294	1307	
01052=4	CIZP	회가 하면 있습니다. 아니라 살이는 것이 그리 경험을 주었다. 결과
0105294	1247116	이 아이트를 하고 하고 있는 말이 살아 있다. 그런 중에는 이번 걸로 보고 있다.
0105294	1237CB	이 보냈다고 하는 말을 하지 않는 얼마 하게 모일이 되었다고 하고요?
0105294	HCB	즐거지 않는데 나를 하고 있다면서 얼마를 보는데 없다.
0105274	10X	하는 여러 아무는 그들은데 얼마라 호른 나는 얼마를 하나 다른다.
0105294	13000	
0105194	incon	
0105294	11121	그는 그리는 살이 보고 있는 생활이 한다면 하는데 아름다고 하는다면
0105254	Frace *	사용하다 마다들이 안 되었다. 물로를 내려가는 된 이번 회사를 모르는
0105274	opca	
0105774	TEB	왕마를 하는 경기 하다 하나 생님, 맛있다. 하나 다쳤네. 아이를 하다 하는 것
0105274		로 공유를 보다 되었다. 그는 그런 그릇이 모르게 되었다.
0105294	IPB	불교하이는 그는 발표를 이 있었다. [2012년 1일 2012년 1
0105270	ocr	경기들 수면 있다는 이번 경우를 살아가면 하다라고 있었다.
0105294	FČT	
0105294	SBC	에서 보고 있습니다. 그런 그렇게 작가 나를 통해를 하는 것이다. 하지만 하는 것들
0105294	DCF:n	
0105294	iloup	회사 전환 (경기) 사람이 되었습니다. 그렇게 함께 모든 하는데
0105294	iicp	경기가는 시간 경기를 걸고 빨리고 하다 되기, 중요한 경송
0105294	1170	
0105294	DEM	보다 경향 전 소전 하지만 수 있다면서 사람들은 하는 것이다.
The second secon	MANUAL CONTRACTOR	이번 회의 모든 의학 에 이탈 회사 속의 불의 출범 모습이 없어만 됐

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O105295 ODCD O105295 SDC8 O105295 SDC8 O105295 120CE O105295 BDC9 O105295 BDC9 O105295 CDB O10	The state of the s		Fee					
0105295	"我,你就想要什么,我,基本一起了一点?	nres						
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0105275 BDCR 0105295 12DCF 0105295 CMSC3 0105295 PM 0105295 CB 0105295 CB 0105295 VC 0105295 VC 0105295 DEM 0105295 DEM 0105295 13UCP 0105295 DEM 0105295 TE 200.00 0105295 TE		the first of the control of the cont						
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0105295 TTE 200.00 0105295 TTEDC 0105295 CTPU 0105295 TEDCP 0105295 PX 0105295 PX 0105295 PN 0105295 PN 0105295 POR 0105295 CDBG 0105295 CDBG					l Pari			
0105295			200.00					
0105275								
0105275 1207F 0105275 11DCP 0105275 PX 0105275 129776 0105275 ECT 0105275 ECT 0105275 ECT 0105275 CDBH		and the second of the second o						
0105275		1227						
0105295		11DCF			in a second			
0105275 124176 0105275 FCH 0105275 FCH 0105275 FCFM 0105275 CD8F: 0105275 TCL	0105075							
0105295 ECH 0105225 HBR 010525 DCFR 010525 CDBH 0105255 TCL								1.50 1.50
0105225 RB8 6165255 DCFR 0105255 DDBH 0105255 TCL								
C165I95 D105295 CD8H D105295 TCL		and the second s						
0105295 CD684								
사람이 사 리 6:2.75 - 이 전 TC L - 이사 있는 하고 그 전략 하나는 이 하는 사람들은 전략이 모든 사람들은	C165295	1. April 10 Miles						
	0105295						57 57 57 1 45 7 5	
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수도 하는데 그 일 사람이 있는 것이 말 중요. 이번 그는 일 사람들은 그는 사람들은 이 그는 사람들은 사람들이 가는 수 있다.								
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injoice No.: 015:1194 Our Lab # 3 0105276 Our Lab W Your Sample 1): Hidr Sample Hatrix : SOIL 1 popular Date: 05/30/90

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\$ 200.00

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T . . :

Attes

Our Lab # : 0105297 Your Sample ID: HIVAY Sample Matrix : 50th Involce 16. : 0103195

Invoice bate: 05/30/90

Fes

\$ 200.00

TOTAL \$ 200,00

ITEMIZED JEST CHARGES: Lab# Yest

> 0105297 11DCE 0105297 HDCE 0105297 PTCE

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0105297 IPU 0105297 I2DCE 0105297 1112TC

0105297 CHERT 0105297 TCE 200-00 0105797 i12T(E

0105297 1231CN 0105297 11220E 0105297 3CLE

0105297 123:F 0105297 HDCH 0105297 HPF 0105297 VC

0105297 BM 0105297 CB 0105297 FTCM

0105297 DCFT 0105297 13CF 0105297 TOL

0105197 65 0105297 CHCLJ 0105297 CH 0105297 SEB

Lab#

0105097

Fee

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0105277	역 (: : : : : : : : : : : : : : : : : :
0105257	200 <mark>1-291-48</mark> -2-19-25-20-25-20-25-20-20-20-20-20-20-20-20-20-20-20-20-20-
0105297	시 (Yet) (^) 사람이 있는데 말라고 하는 하는지 않는 나를 되었다.
0105297	- 1 . in : 1
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010527	크리 주마 열리는 전문하실 수 있었다. 이 모를 수 없는 물론에
0105297	보고 (2) (b) () : [[[[[]
0105297	님은 1700년 다른 사람들은 그는 얼마스 사람들은 말을 가는 사람들이
0105297	
0105297	
0105257	다 EP 이렇게 다음하게 된 하셨습니다. 스탠딩을
0105297	H - ST 하는 이 보고가 있는 그런, 항상들하지 않았다. 함께 1
0105297	- A. Tuz: c - B B B B B B B
0105227	
0105297	2 (1 1.10) [15] - 보는 시간 (12.20] (1.20] (1.20) (1.20)
0105297	트리스 트웨 트 : 11 10 10 10 10 10 10 10 10 10 10 10 10
010578	
0105297	하는 발생이 말이라면 하는 않아 보았다. 그렇게 살아 있다.
0105297	(BELORE) : 10 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1
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0105297	변나CCT 이 전 Back Park Park I I I I I I I I I I I I I I I I I I I
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0105297	문을 22PGP 이 사용 문을 넘기 보이는 일을 넘기 들었다. 참시가 취용된

To:

Atto:

TOTAL

\$ 200.00

ITEMIZED TEST CHARGES: F. ... 12411 E 0105298 T1200 0105298 CCIT 0105298 LIDEP 0105293 11CD 0105298 130P 0105298 ODCE 0105298 111TCE 0.20 0105298 VC -0105298 TCE 200.00 0103298 Ph 0105278 CHBFIS 0105299 OX. 0105298 PDCB 0105273 12rcF 0105258 220CP 0105293 EB 0105258 0126 0105278 CIT 0105299 CT 0105278 120CE 0105298 FEZ 0105293 DBH 0105298. HOCE 0105298 124TOR 0105298 123TCB 0105293 (i.f.) 0105299 EOCH 0105278 F'X 0105298

Laboration of the contract of

0105298

TO:..

		그림하다 내가 가지 않는데 이상 생생이 있다. 스토스웨티 나는 사실 입사를 보냈다.
Alaba in an an a	Tost	
بيباء الشبك بالباب والساعة سنستشب		
0105293	. # (Y * * * * * * * * * * * * * * * * * *	마음 등이 돼지가 이번 그렇게, 전화 등을 보고 말을 보고 되었다. 중
6105296		하마스트로 하면 있는데 나는 중에 그릇을 보는 것 같아. 그들까 이 모델
0:05255	COST	
6105298		하람들이 그 젊으는 회원 등 나가장이 많아 되는 것 같아 해가겠습니다.
3927010	CHELLA	이 경기에서 가는 눈이 들어 못하는 말로만 먹고하다면 다른데
0105273	CCF 11	
0105298	HCE	보기를 통해 하는 사람이 되었다. 나는 사람들은 살아 없는 사람들은 얼마를
0105278	are .	보고 본 하는데 말했다면 하는 네트를 받아갔다면 없는 말을 제 등으로 하는데
0105298	11227E	하는 회장 학교를 통해 고면 되어 놓으면 다녔으면 하는 일시 회사들이 모든
0105298	1 3511ec	교레 라마레 아시아는 아이를 다른 경우를 하는 것 같은데 모든 모든 모든 것
0105298	1127CE	
0105278	14100 12 44	일본인 중점이 되었다. 그들은 하늘 만큼 하는 이, 하는데 하나 하는
0105298	ECIT	힘들 하고 어디를 했다면 하늘에 다일까지 않았다. 이 과학 등 년
0105298	116.5	
0105298	OCLE .	이 아마이지 않아야 되는데 그래 하다 그들은 것 같아요?
0105298	123TP	물 그렇다 속한 다른 살을 살으면 그리면 없다. 그렇다면 했다는 그림
0105298	IBE	요시 그리다 사람이 가지만 사계를 하는 것으로 하게 되고 않는
0105298	FCT	
0105298	CB	
0105298	MX (경영 하는 바다 하는 사람들은 사람들은 바다 하는 것이 없었다.
0105298	# \$ 2000 174	네 공기를 가지 않는데 되는 소개 있다. 하라고 함께를 게 되었다. 네 요즘
0105298	, 1117.VE	
0105298	Fif B	[일을 하는 회송] [2011년 10년 12일 22일 22일 22일 22일 22일 22일
0105298	FICH	사용하는 이글 사용하는 사용 중요 맛을 되는 글로운 사용을 했다.
0105298	OCT	이 나는 뭐 그는 것이 가지를 하지 않는 것들이 하는 것이 하는데 이 것은데 없다.
0105298	SER	그 회사님이 되는 이번 않는 한 경험 보이고 있는 것이다.
0105298	PIST	보이 동생님이 작은 마음에서 그래 있다. 경기하지 않다고 말이 됐다. 전기주 시간 연

- THOOSEE

To:

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IHVOICE SUMMARY:

L.ab#

0105279

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ITEMIZED MEST CHARGES:

£ 01:05227 Our Lab H Your Sample To: HIWA

Sample Waters : SOIL

Test Charges 429.00

 $\Gamma \oplus \Xi$

Additional Charges:

429.C

429 TOTEL

0105299 NH IF TH 0105299 DEF HHOPA 0105299 F ,-, 0105299 0105299 HaE 0105297 HOB HCE 0105299 246647 0105207 .0. 0105299 FIFTH . 4CFFT 0105279 0105200 Act 11 F 7 0105299 1247CE 0105299 1 0105299 A#E 185.00 0105299 BEHP 0105299 0105299 DHOF ;=·;=. 0105299 0105299 4:47 0105299 BOE 1000 B 0105299 0105279 FIR 12100 0105299 0105297 BCIPE

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Invoice No.: 0103197

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0105259	2CH
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0105299	CdE

LAVOICE

To:

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0105300 ECE 225.00 T 0105300 0105300 120CE 17:-0105300 0105300 FCEM 140CB 0105300 0105300 DEA 0105300 EUFA 0105300 BID 0105300 13000 2378 D 0105300 0105300 4CPPE 0105300 BKFA 0105300 ro. 124TCB 0105300 0105300 FA 0105300 ACH 0105300 Dispu 0100300 E.E.F. 0105300 Fine 0105300 CEY $S \pm \Xi$ 0105300 0105300 1 :E: HEE 0105300 0105300 NNETA 0165300 120PH 0.05300 AFFEE 0105300 HCEZ 3105300 24111

HOB

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0105300	ANT					
0105300	330CB					
0105300	DPA					
	 		물론 경제 없			

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· THVOTCE

Attn:

!m/wice Hc.: J103192 Our Lab # : 0105301 Your Sample ID: HTWAY In sice Date: 05/30/9 Sample Batrix : 501L INVOICE SUMMARY: Test Charge: 429.00 . Additional Charges:

NIZED TEST	CHAROES:	Fee					
0105301	aka c E						
010579							
01:5301							
0105301	HEER						
0102201	DE F						
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0103301	4CPPE						
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Ojostoj	Correspond
Oleszár	
0105301	회 리: (() 이 나는 그림, 작물하는 사람들은 사람 모양하다
6105301	의 성에 에 보인트등 없는 항상의 사람들이 등 선택 등 등 있다.
0105701	
1 - 1 - 1 - 1 - 1 - 1 - 1 - 1 - 1 - 1 -	
0105301	
0105301	BAPE 이 경기에 진유된다고 말하다고만 하다.
0105301	COLUB CE COLOR COLOR COLOR COLOR COLOR COLOR CE COLOR CO
0105301	고 보는 100 등 일본하고 말을 하는 것을 다시 수 있다.
0105301	December 1925. Out the Ring of the Fig.
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0105301 0105301	
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0105301	된 # 호텔 : 호텔 : 살림 : : : : : : : : : : : : : : : : :
010530;	'비HCE' 등 등 등 하는 사람들은 사람들은 바람들이 되었다.
0105301	TROIPE

- IN OTE E

To:

Attria

Our Lab # : 01/5302 Your Sample ID: HTWAY Sample Matrix : 50 L Invoice Date: 05/30/90

TIEMIZED TEST CHARGES:

Lab# Ye#1 Feo

0105302 **F**. 0105302 12008 0105302 CJE 0105302 HCWZ 0105302 CrE 0105302 BEHR 0105302 FA 0105302 140CB 0105302 PAG 0105302 CRT 0105302 SIDCE 0105302 OBFT 0105302 HNMA 0105302 FA 6105302 FIT 0105302 SEFFE 0105302 72DFH 0105302 237970 0105302 AgE 0105302 HaE 0105502 FE. 19.00 0105302 Addit 165.00 0105302 ESE 0105302 130CB 0105302 E.BF 0105302 BAL

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0105302 FME 0105302 FME 0105302 CMF 0105302 CMF 0105302 FLP 0105302 FLP 0105302 DNOF 0105302 BEFA 0105302 HMPM 0105302 I 0105302 HMPM 0105302 RCH	1	D105202	HOT	
O105302 FME O105302 SMF O105302 DMF O105302 FLP O105302 FLP O105302 DMGP O105302 BEFA O105302 BEFA O105302 HMSPR O105302 H HCE O105302 HCE O105302 BCEM O105302 B		0105302	en. Cr	[설송이는 12년째] EN 20년 등 12년
0105302 94E 0105302 94F 0105302 FLB 0105302 FLB 0105302 BBFA 0105302 BCE 775.00 0105302 HNPM 0105302 I 0105302 HCE 0105302 RCEM 0105302 RCEM 0105302 BCFB	-	0105702	firf a	경을 하늘의 이번 내는 것이 맛있다. 하지 않는 그런 이번에 모르겠다면
0105302	:.	0105392	非成 点。	요즘 좋은 하면 보고 하다 하는 경험이 얼굴하는 것 같아요?
0105302 250HT 0105302 FLF 0105302 DMOF 0105307 BEFA 0105302 BCE 225.00 0105302 HCE 0105302 HCE 0105302 BCEM 0105302 BCEM 0105302 BCEM 0105302 BCIFE		0105302	SeE	고등을 존속하실 때문에 다려면 다 열심을 다리가 하다고 말을
0105302 FLP 0105307 BEFA 0105307 BEFA 0105302 BCE 278.00 0105302 1 0105302 HCE 0105302 RCFH 0105302 BCIPE 0105302 BCIPE 0105302 SHIP 0105302 SHP 0105302 BCIPE 0105302 BCIPE 0105302 BCIPE 0105302 BCIPE 0105302 BCIPE 0105302 BCIPE	, ;	0105302	CONTRACTOR OF STREET	존리하다 하다 가는 내용에 가장 살아서 모든 사람이는 것이 없다면서
010530? DNOF 010530? BEFA 010530? HEFT 010530? ACH 010530? HCE 010530? BCEM 0105302 BCEM 0105302 BCIFE 0105302 BCIFE 0105302 BCIFE 0105302 BCIFE 0105302 BCIFE 0105302 BCIFE 0105302 BCIFE 0105302 BCIFE 0105302 BCIFE 0105302 BCIFE		0105302	2500	그렇게 된 그만들다 아름일이라고 된다. 나는 끝 나다
0105307 BBFA 0105302 BCE 025.00 0105302 I 0105302 HCE 0105302 RCEM 0105302 RCEM 0105302 BCIPE 0105302 BLD 0105302 BLD 0105302 BLD 0105302 BLD 0105302 BLD 0105302 BEP 0105302 BEP 0105302 BEP 0105302 BEP	÷.	0105302	FLF	
0105302 PCE 275.00 0105302 PMPM 0105302 PCH 0105302 PCEH 0105302 PCEH 0105302 PCEH 0105302 PCIPE 0105302 PCIPE 0105302 PCH		0105302	DHOF	결과에 그리고 하고 있다. 그는 그리는 말을 걸 때 없는데 그 없다.
0105302 HEPE 0105302 I 0105302 HCE 0105302 RCEM 0105302 RCEM 0105302 RCIPE 0105302 RCIPE 0105302 RIG 0105302 RIG		0105305	BEFA	생활 중에 하는 하고 있는데 나로를 들는 점을 살으면 하여 모모를 했다.
0105302 1 0105302 #CH 0105302 #CEH 0105302 :24TCE 0105302 %CIPE 0105302 24DNT 0105302 & RIG		0105365	ECE 270	나 아 이번 시간에 있는데, 그를 들어보는 말은 사람이 살아 있는 말을 받는다.
010530? ACH 0105302 HCE 0105302 ;24TCE 0105302 WCIPE 0105302 24DHT 0105302 kiG 0105302 ECH 0105302 DEP 0105302 DEP 0105302 HMOPA 0105302 HMOPA		0105302	HMPH	
0105302 HCE 0105302 RCEM 0105302 SCIPE 0105302 RAIG 0105302 RIG 0105302 RIG 0105302 RIG 0105302 REM 0105302 REM 0105302 REM 0105302 REM 0105302 REM		0105302		그렇게 하를 가장하는 다음을 가지만 못 못하면 적인을 아닌다면?
0105302 RCEM 0105302 SCIPE 0105302 SAINT 0105302 RIG 0105302 RIG 0105302 DEM 0105302 DEM 0105302 HNOPA 0105302 HNOPA		0105302	#CH	하면 선생님이 이렇게 되었다. 이렇게 하면 하게 되었다.
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WATER WELL SURVEY

WELL INVENTORY ALABAMA AIR NATIONAL GUARD FACILITY DANNELLY FIELD MUNICIPAL AIRPORT MONTGOMERY, ALABAMA

Prepared for

AIR NATIONAL GUARD READINESS CENTER ANDREWS AIR FORCE BASE, MARYLAND

Submitted by

CH2M HILL, INC. 2567 FAIRLANE DRIVE MONTGOMERY, ALABAMA

July 1995

Introduction

As part of the Site Investigation for the Alabama Air National Guard (ANG) at Dannelly Field in Montgomery, Alabama, a well inventory was conducted. The purpose of the well inventory is to identify potential receptors that may potentially be affected by compounds that may be identified during the Site Investigation.

The inventory was conducted for an area extending 1 mile out in all directions from the ANG Dannelly Field Facility. The estimated aerial extent of the well inventory is shown graphically in Figure 1. In addition, all public water supply wells located within a 5-mile radius at the ANG were identified.

Methodology

Initially, the study area was identified on a United States Geological Survey (USGS) 7-1/2 minute topographic quadrangle. This map was used to identify the well inventory area and as an aid to locate households and businesses currently using water not supplied by the City of Montgomery.

Before the door-to-door survey, a plat of the area's municipal water supply systems, supplied by the Montgomery Water Board, was reviewed. To supplement the review process, a windshield survey was conducted within the 1-mile radius of the study area. Both recent and more established developments were identified, looking for pump houses and other evidence of private water usage. On the basis of the plat and the assumption that the established developments are more likely to use private water, they were targeted for the door-to-door survey.

The door-to-door survey was conducted at the households identified during the windshield survey. The interviewers visited 43 households within the area shown in Figure 1. The interviewers repeatedly tried to talk to several residents, but some were not at home during the survey. However, the survey team was able to establish their water source through discussions with surrounding neighbors.

Results

The water use in the area is a combination of municipal water systems and privately owned wells, with municipal water being the primary source. The private wells identified are installed in either the Eutaw aquifer (deep well) or in shallow sand and gravel beds located within the terrace deposits (shallow well). Many of the respondents were unsure

of the actual depth of their wells, thus making it difficult to obtain an exact number of deep and shallow wells located within the area. Additionally, several residents indicated that they obtained their water supply from naturally occurring springs and artesian wells located within the study area.

The results of the well inventory are summarized below:

Operating wells	17
Abandoned wells	15
Naturally occurring springs and artisan wells (used as a water source)	2

Table 1 lists the summary of the door-to-door survey. Figure 1 shows the estimated aerial extent of the well inventory area as well as the location of the operating wells, abandoned wells, and springs. Each identified operating well and spring delineated in Table 1 was assigned a number corresponding to the numbered locations shown in Figure 1. The operating wells are identified in Figure 1 by circles, abandoned wells as triangles, and the spring (water supply) as a square. Figure 2 shows the location of public water supply wells operated by Montgomery Water Board.

Discussion

Two main areas are shown on the east and northeast sides of the ANG as being served by privately owned water sources (Figure 1). One well was also identified southwest of the ANG, but it is solely used for watering livestock (Well 19, Figure 1). The City of Montgomery is currently extending its water service in both of the private water usage areas. Residents are or will be given the opportunity to switch over to municipal water service. Furthermore, the Montgomery Airport Authority is actively purchasing property located within the 1-mile radius and tearing down the existing dwellings, thus reducing the number of private wells. This purchasing, coupled with the increasing availability of municipal water, will mean a continued reduction of private water users within the 1-mile radius.

The area shown to the northeast of the ANG along Old Hayneville, Brewer, and Rich Roads shows a concentration of private water users. The wells found in this area are screened at various depths, reportedly ranging from 40 ft to more than 200 ft. Additionally, a spring and a well under artesian conditions were identified in this area.

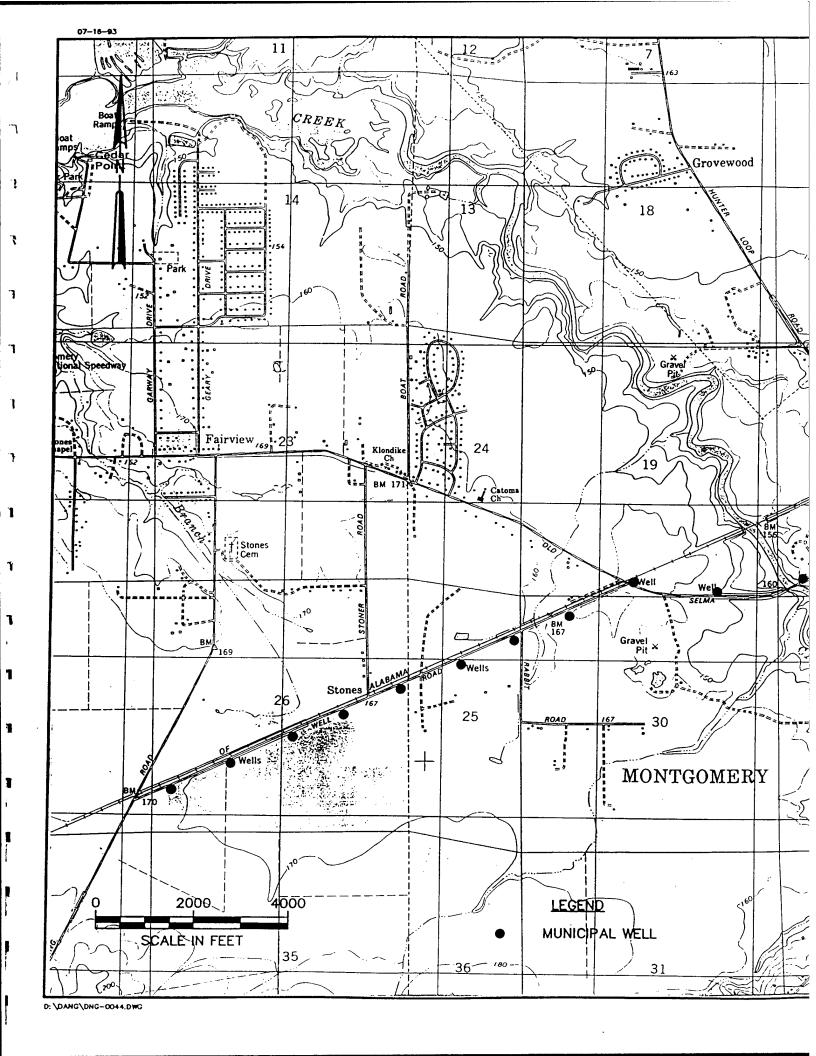
Table 1 Dannelly Field, Montgomery, Alabama Source Summary

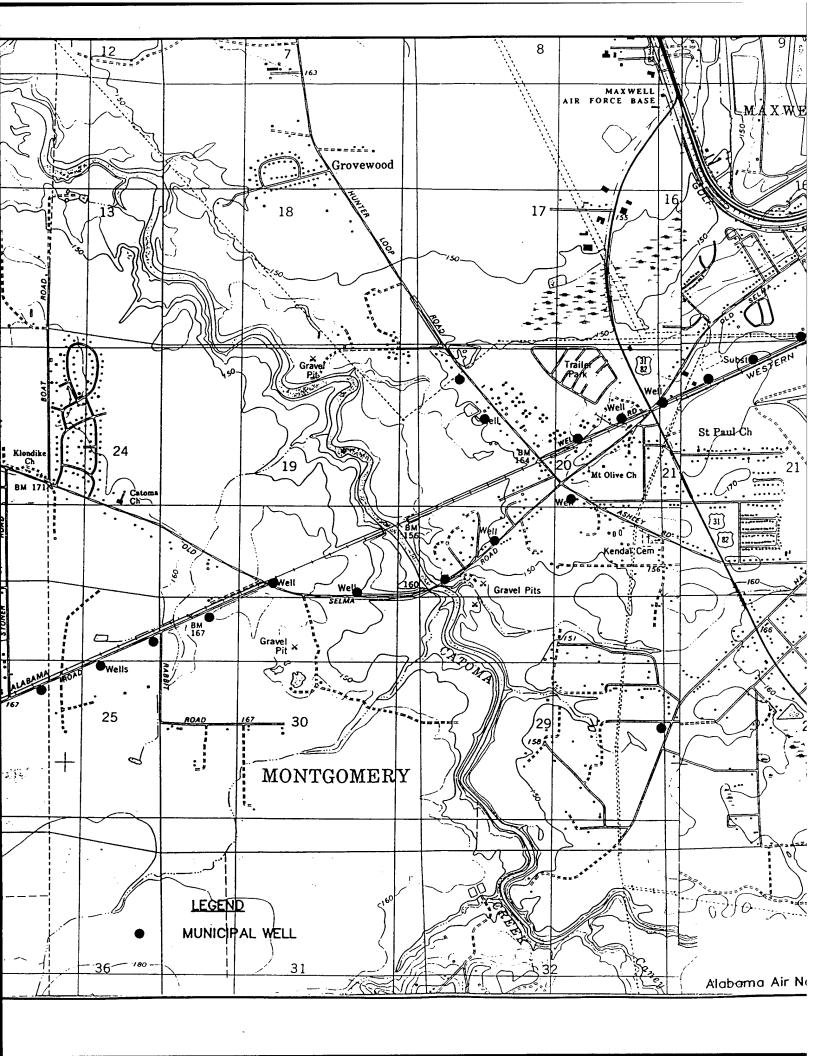
Identification				Approximate
Number	User Address	Source	Use	Depth, (ft)
1	4620 Old Hayneville Rd	spring	consumption/irrigation	ground surface
2	4560 Old Hayneville Rd	well	consumption/irrigation	>200
3	3915 Brewer Rd	well	consumption/irrigation	200
4	4328 Brewer Rd	well	irrigation	80
5	4340 Brewer Rd	weli	consumption/irrigation	>150
6	4412 Brewer Rd	well	consumption/irrigation	unknown
7	4420 Brewer Rd	well	consumption/irrigation	40
8	4458 Rich Rd	well	consumption/irrigation	>150
9	4460 Rich Rd	well	consumption/irrigation	unknown
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	4512 Rich Rd			
11	4524 Rich Rd	well	consumption/irrigation	85
12	4530 Rich Rd	well	consumption/irrigation	>200
13	4561 Rich Rd	well	consumption/irrigation	unknown
14	4554 Rich Rd	well ^b	consumption/irrigation	unknown
15	1250 Old Lamar Rd	well	consumption/irrigation	108
16	1180 Old Lamar Rd	well	consumption/irrigation	unknown
17	1185 Old Lamar Rd	well	consumption/irrigation	65
18	1145 Old Lamar Rd	well	consumption/irrigation	180
	Rt. #1, Box 22B			
19	Hope Hull, AL 36043	well	watering livestock	unknown
•	(334) 286-0931°			

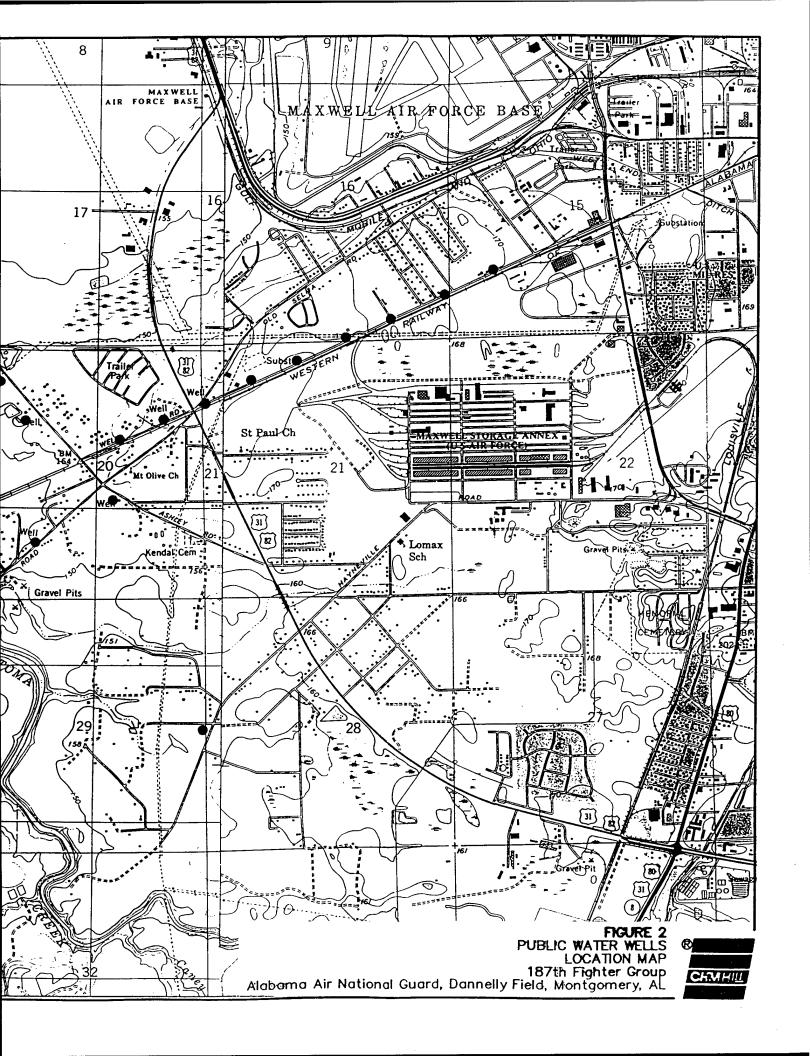
^aWell serves both residences

 $b_{\mbox{Well under artesian conditions}}$

^COwner's contact address and phone number







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The area to the east of the ANG also shows a concentration of privately owned wells. The wells in this area (Old Lamar Road) are screened at various depths and reportedly range from 65 ft to 180 ft. However, a number of the residences did not know the depth of their wells. The wells in this area probably are screened within the Eutaw Formation below the Mooreville chalk.

All of the private well (deep and shallow) and spring water users were asked to comment on the quality of their water. They all considered the water to be superior to the municipal source and noted a reluctance to switch. There has been no apparent decline in the water quality over years of consumption.

The remaining areas within a 1-mile radius of the site are served by a public municipal water system. The interviewers visited a few households in each of these areas and learned that the areas are served by these water systems. Therefore, they did not visit the remaining households in each area.

Within a 5-mile radius of the ANG installation, the Water Works and Sanitary Sewer Board of the City of Montgomery operates the West Well Field. The wellfield is located primarily along Foshee and Well Roads at a distance ranging from 2 to 5 miles north of the ANG. The wellfield (Figure 2) includes 29 wells that currently supply about 36 percent or 21 million gallons per day of potable water for the Board. The wells are typically multiple aquifer completions that withdraw water from the Eutaw, Gordo, and Coker Formations at an average depth of approximately 700 ft.

The Water Board supplies about 5 percent of the remaining balance of its water supply from the North Well Field and about 59 percent from the C. T. Perry Plant, which treats water from the Tallapoosa River. The North Well Field and the Perry Plant's raw water sources are located in the northern and northeastern Montgomery areas, respectively.

TOXICITY VALUES OF SELECTED CONTAMINANTS

APPENDIX I Toxicity Profiles for Human Health Effects for Chemicals or Potential Concern Alabama ANG--Dannelly Field

ACETONE

Acute exposure to acetone can cause irritation of the eyes, skin, and respiratory tract. Depression of the central nervous system and narcosis can occur after inhalation of high concentrations of acetone. Dryness of the mouth and throat, dizziness, nausea, incoordination, loss of speech, and even coma have been described in some cases of workers exposed occupationally to acetone. In a chronic study reported by EPA (IRIS, 3/1/88), kidney damage was observed in albino rats administered acetone by oral gavage. Kidney damage and metabolic changes have been noted in humans who ingested acetone (Sax, 1989). Acetone can increase the toxicity of other chemicals, particularly solvents. The hepatotoxicity of carbon tetrachloride can be increased greatly in the presence of acetone.

ANTIMONY

Many antimony compounds irritate the gastrointestinal tract. Acute intoxication with antimony results in severe vomiting and diarrhea. With occupational and inhalation exposure, rhinitis and acute pulmonary edema may occur. The chronic inhalation of some antimony compounds can produce respiratory problems and diseases. Transient spots on the skin have been reported in some workers. Antimony may form stibine gas, which causes hemolysis.

ARSENIC

Acute oral exposure to arsenic can cause muscle cramps, facial swelling, cardiovascular reactions, severe gastrointestinal damage, and vascular collapse leading to death. Sensory loss and hematopoietic symptoms delayed after exposure to high concentrations are usually reversible. Inhalation exposures can cause severe irritation of the nasal lining, larynx, and bronchi. Chronic oral or inhalation exposure can produce changes in skin, including hyperpigmentation and hyperkeratosis; peripheral neuropathy; liver injury; cardiovascular disorders; oral exposures associated with peripheral vascular disease; and blackfoot disease. Arsenic is a known human carcinogen; oral exposures are associated with skin cancer, inhalation exposures with lung cancer. Toxicity varies for different compounds; inorganic trivalent arsenic compounds are usually more toxic than pentavalent compounds; high doses of some inorganic arsenic compounds to pregnant laboratory animals produced malformations in offspring.

BARIUM

Acute ingestion of barium salts can cause prolonged muscular stimulation, gastroenteritis, hypokalemia, and cardiovascular effects such as ventricular fibrillation and extra systoles. Prolonged occupational inhalation has resulted in baritosis-a benign, reversible pneumoconiosis. Toxicity of compounds depends on solubility.

BENZENE

Acute exposure to high concentrations of benzene vapors may cause central nervous system depression, unconsciousness, and fatal cardiac arrhythmias. Benzene is readily absorbed after oral and inhalation exposure. Chronic exposure of workers to low concentrations has been associated with blood disorders. Chromosomal aberrations in bone marrow and blood have been reported in experimental animals and some workers. Benzene has been classified as a Group A Human Carcinogen, based on evidence of carcinogenicity from epidemiological studies and correlations between exposure by inhalation and leukemia.

BENZO[A]ANTHRACENE

No information is available on short-term dermal or inhalation effects. There is no information on systemic effects. PAHs as a group may cause skin disorders and immunosuppressive effects. Evidence exists that benzo[a]anthracene is carcinogenic to laboratory animals through dermal and ingestion exposure routes. Inhalation data are not available. Benzo[a]anthracene may cause skin and lung cancer. No reports exist relating cancer in humans from exposure to benzo[a]anthracene exclusively, but there are reports on exposure from PAH mixtures. This chemical is mutagenic in laboratory experiments. Benzo[a]anthracene may be metabolized into reactive derivatives.

BENZO[A]PYRENE

Acute toxicity for Benzo[a]pyrene appears low when administered by oral or dermal routes to laboratory animals. Prolonged exposure may produce chronic dermatitis and reproductive changes. Repeated oral doses to mice have caused hypoplastic anemia. Induction of cancer is the key toxic endpoint from intermediate and long-term exposure. Benzo[a]pyrene is a constituent of coal tar, which is classified as a Level 1 known carcinogen by IARC and a Level B2 probable carcinogen by the EPA. Ingestion may produce stomach tumors, and inhalation may produce lung cancer. Prolonged skin exposure has been linked to an increase in skin cancer among workers. Benzo[a]pyrene is considered to be the most potent carcinogenic PAH. Benzo[a]pyrene is a mutagen.

BENZO[B]FLUORANTHENE

No acute toxicity information is available. Systemic effects specific to benzo[b]fluoranthene have not been reported. Skin disorders and immunosuppressive effects have been reported for PAH mixtures. Experimental evidence exists that it causes lung and skin cancer in laboratory animals by dermal absorption and intratracheal distillation. There is no evidence of reproductive or teratogenic effects.

BENZO[G,H,I]PERYLENE

Limited information for acute or chronic toxicity is available. It is a liver and skin carcinogen in laboratory animals. Data available are inadequate to determine carcinogenic potential in humans. Benzo[G,H,I]perylene has been reported to produce cocarcinogenic effects when applied to mouse skin along with benzo[a]pyrene.

BIS(2-ETHYLHEXYL)PHTHALATE

(Di(2-ethylhexyl)phthalate or DEHP or BEHP)

In general, low acute toxicity has been reported in experimental animals; accidental acute exposure in man resulted in mild gastric disturbance and catharsis. Chronic exposure at relatively high concentration has retarded growth and resulted in increased liver and kidney weight in experimental animals. Oral administration to rats and mice resulted in increased hepatocellular carcinomas or neoplastic nodules. Classified by EPA as a B2 carcinogen. Some evidence in animals of teratogenic and fetotoxic effects exists. Reproductive effects, decreased fertility and testicular damage have been noted in rodents. This substance is poorly absorbed through skin and is rapidly metabolized.

BROMODICHLOROMETHANE

Bromodichloromethane results from chlorination of precursors in raw water. Toxic effects of bromodichloromethane have been shown in laboratory animals and include sedation and hemorrhage in the kidney, adrenals, lungs, and brain. It is classified as a Group B2 Probable Human Carcinogen by ingestion exposure routes.

BROMOFORM

Acute exposure to high levels of vapor produces irritation of the respiratory tract, pharynx, and larynx, with lacrimation and salivation. Other effects include headache,

listlessness, vertigo, unconsciousness, loss of reflexes, convulsions, and death usually caused by respiratory failure. This substance can cause decreased liver function and histopathological effects in experimental animals. Longer-term exposure to mice by oral administration causes decreased body weight, fatty metamorphosis of the liver, and hepatocellular changes. Inhalation studies with rats show disorders in prothrombin synthesis and glycogenesis in the liver and reduced filtration capacity in the kidney. Bromoform results from chlorination of natural organic precursors in raw water.

CADMIUM

For acute exposures by ingestion, symptoms of cadmium toxicity include nausea, vomiting, diarrhea, muscular cramps, drop in blood pressure, vertigo, loss of consciousness, and possible collapse. Exposure by inhalation can cause respiratory ailments, acute chemical pneumonitis, and pulmonary edema. The major chronic effects in workers are respiratory and renal toxicity. Cadmium bioaccumulates in the kidney, and nephropathy results after critical concentration (around 200 ug/g) is reached. Inhalation can cause chronic obstructive pulmonary diseases. Chronic exposure also affects calcium metabolism. Cadmium is classified as a Group B1 Probable Human Carcinogen, by inhalation routes.

CHLOROFORM

Acute toxic exposure effects of chloroform in humans include respiratory depression, coma, and liver and kidney damage. Chloroform (anesthetic) depresses the central nervous system and may result in cardiac arrest (apparently from sensitization to epinephrine). The chronic toxic exposure effects in animals include liver and kidney damage, fetotoxicity, and malignant tumors.

CHROMIUM

Chromium is an essential micronutrient and is not toxic in trace quantities. Following oral exposure, absorption of chromium (III) is low while absorption of chromium (VI) is high. The major acute effect from oral exposure is renal tubular necrosis. Inhalation of chromate salts can cause irritation, inflammation, and ulceration of nasal mucosa. Chronic exposure to chromium (VI) can result in kidney damage. Inhalation exposures in industrial settings have resulted in damage to the respiratory system and have been associated with excess lung cancers. Exposures to the skin can result in allergic skin reactions in sensitive individuals. Overall, the hexavalent forms are usually more toxic than trivalent forms. Inhaled chromium (VI) is classified in Group A as a human carcinogen.

CHRYSENE

Chrysene is absorbed by oral and dermal doses and accumulates in adipose and mammary tissues. Chronic toxic effects have not been described. It is carcinogenic in laboratory animals exposed to long-term dermal doses. There is limited evidence that chrysene is mutagenic. Epidemiological reports document incidences of skin cancer when exposed to PAH mixtures that included chrysene.

COPPER

Copper is an essential micronutrient and is not toxic in quantities necessary for human health. Acute inhalation exposure to copper dusts can result in symptoms similar to metal fume fever. Exposure to dusts and mists of copper salt results in nasal congestion. Exposure to fumes results in upper respiratory tract irritation, metallic or sweet taste, and skin and hair discoloration. The major chronic toxic effect is hemolytic anemia in some dialysis patients.

DIBENZO[A,H]ANTHRACENE

Oral absorption occurs but there is slow dermal absorption. IARC [1983] has concluded that there is enough evidence that dibenzo[a,h]anthracene is carcinogenic to laboratory animals. In laboratory experiments, oral doses have caused tumors in mice; intratracheal distillation has caused lung tumors in rats and dermal application has caused skin cancer. High doses in laboratory animals have produced fetal deaths.

DIBROMOCHLOROMETHANE

In laboratory animals, dibromochloromethane causes liver changes including fatty metamorphosis, calcification, centrilobular necrosis, and vacuolar changes. Other effects include toxic nephropathy, salivary gland inflammation, and thyroid hyperplasia. It results from chlorination of precursors in raw water. There is some evidence of mutagenicity.

DI-n-BUTYL PHTHALATE

One case of human oral exposure has been reported to cause nausea, vomiting, dizziness, headache, pain and irritation in the eyes, lacrimation, photophobia, and conjunctivitis; there was some renal involvement followed by recovery within 2 weeks. Irritation of eyes and upper respiratory tract has been noted in mice that inhaled aerosols of the compound. Liver and kidney lesions have been observed in mice receiving chronic oral doses. No carcinogenicity data are available. Di-n-butyl

phthalate may cause increased embryotoxicity and teratogenic effects in rats and mice. Testicular atrophy observed in animals may result from disturbances in zinc metabolism.

1,1-DICHLOROETHANE

1,1-Dichloroethane may cause central nervous system depression when inhaled at high concentrations. It is also a skin irritant. It has been classified as a Group C possible human carcinogen by ingestion exposure routes.

1,2-DICHLOROETHANE

CNS depression, lung irritation, and injury to liver, kidney, and adrenals have been reported. Deaths in humans exposed by ingestion or inhalation may result from circulatory and respiratory failure. Chronic exposure can cause liver degeneration and kidney damage in laboratory animals. Eye damage (necrosis of corneal epithelium) has been observed in dogs injected with 1,2-dichloroethane. Repeated exposures have been associated with anorexia, nausea, liver and kidney dysfunction, and neurological disorders in workers. It is carcinogenic in mice and rats exposed orally, and mutagenic in some tests in bacteria, barley, and fruit flies.

1,1-DICHLOROETHENE

1,1-Dichloroethene (Vinylidene Chloride) is rapidly absorbed following oral and inhalation exposure. The liver appears to be the principal target after acute exposures. At high concentrations, inhalation can cause central nervous system depression and unconsciousness. Reports of chronic health effects on workers exposed to 1,1-dichloroethene include liver function abnormalities, neurological and sensory disturbances, weakness, and fatigue. It is described as an "exquisite hepatotoxin" because it is more potent and faster acting than the classic hepatotoxin, carbon tetrachloride. Its structure is similar to vinyl chloride, a known human carcinogen. 1,1-Dichloroethene is classified as a Group C possible human carcinogen.

DICHLOROMETHANE (METHYLENE CHLORIDE)

Dichloromethane acts as a mild narcotic irritating to eyes and upper respiratory passages. Fatalities have been associated with acute or prolonged exposure. In animals chronic exposure can affect the liver and kidney. Dichloromethane may substantially increase carboxyhemoglobin levels, preventing the transfer of oxygen to tissues. Damage to liver and CNS following long-term occupational exposure has been

reported. It is a carcinogen in female rats and male and female mice, is classified as B2 probable human carcinogen by EPA, and is mutagenic in some bacterial tests.

ETHYLBENZENE

Acute inhalation exposure of humans to ethylbenzene concentrations of 435 mg/ml for eight hours has been found to result in sleepiness, fatigue, headache, and mild eye and respiratory irritation.

FLUORANTHENE

Fluoranthene is acutely toxic by oral and dermal absorption. It can cross epithelial membranes and is a defatting agent that may affect the skin. Limited information available. IARC [1983] concluded there is no evidence that fluoranthene is carcinogenic on the basis of available data. When applied to laboratory animal skin simultaneously with other carcinogenic PAHs, it has increased the carcinogenicity of the compound (i.e., cocarcinogenic effects).

FLUORENE

Fluorene may be toxic by inhalation, ingestion, or dermal contact and absorption. Data are inadequate to determine carcinogenic effects [IARC 1983]. No toxicity data are available for humans.

LEAD

Lead is stored in the body in bone, the kidney, and the liver. The major adverse effects in humans caused by lead include alterations in the hematopoietic and nervous systems. The toxic effects are generally related to the concentration of lead in the blood. Blood concentration levels of over 80 ug/dl in children and over 100 ug/dl in sensitive adults can cause severe, irreversible brain damage, encephalopathy, and possible death. Chronic low level exposure to lead can affect the hematopoietic, nervous, and cardiovascular systems. Characteristic effects of chronic lead intoxification include anemia, effects to the immune system, sterility, neonatal mortality, abortion, and morbidity. Children are especially sensitive to low level effects. Acute toxic exposure in humans is characterized by encephalopathy, abdominal pain, hemolysis, liver damage, renal tubular necrosis, seizures, coma, and respiratory arrest. Certain lead salts are classified as Group B2 Carcinogens--Probable Human Carcinogens.

NAPHTHALENE

Inhalation of vapor may cause eye irritation, headache, and confusion. Ingestion may cause abdominal pain, nausea, and vomiting. Skin or eye contact may lead to systemic effects such as bladder irritation, kidney effects, and nemoletic effects such as anemia and decreased hemoglobin. In animal studies, bronchial necrosis was observed in rats. Occurrence of cataracts upon naphthalene vapor and dust exposure has been observed in humans. Subchronic animal studies have shown that oral doses produced cataracts and degeneration of the retina. Dermatitis has been reported with repeated skin exposure. Two studies have reported hemolytic anemia in infants born to women exposed during pregnancy. Studies have not shown that naphthalene is carcinogenic. Naphthalene is commonly found in coal tar and epidemiological studies have shown coal tar to be carcinogenic. The role of naphthalene alone could not be determined. Acute exposures to large doses may cause hemolytic effects (destruction of red blood cells). This effect is most pronounced in individuals with a hereditary deficiency of glucose-6-phosphate dehydrogenase.

NICKEL

Acute exposures to nickel-containing dust may result in chemical pneumonitis. Signs of acute nickel toxicity may include nausea, headaches, vomiting, chest pain, cough, hyperpnea, cyanosis, gastrointestinal and central nervous system effects, weakness, pneumonia, respiratory failure, cerebral edema, and death. Workers chronically exposed to nickel-containing compounds report allergic contact dermatitis and other dermatological effects, rhinitis, and nasal sinusitis and mucosal injury as among the most frequent effects. Nickel compounds implicated as having carcinogenic potential include insoluble dusts of nickel subsulfide and nickel oxides, vapor of nickel carbonyl and soluble aerosols of nickel sulfate, and nickel carbonyl.

PHENANTHRENE

Phenanthrene is an irritant through inhalation and ingestion exposure; it may also be dermally absorbed and an allergen. There is inadequate data for the evaluation of cancer potential in experimental animals. It can cause photosensitization of the skin.

PHENOL

Phenol is corrosive to tissue. Severe eye damage and blindness may result from direct eye contact. Skin contact may produce whitening of skin, burn, or systemic poisoning. Paleness, weakness, sweating, headaches, cyanosis, kidney damage, and death may occur. Chronic phenol poisoning is rare. It induces vomiting, difficulty swallowing,

diarrhea, lack of appetite, headaches, fainting, dizziness, and neural disturbances. Liver and kidney damage may occur. Phenol may promote the effects of certain carcinogens.

PYRENE

Limited information is available. Evidence suggests that pyrene is cocarcinogenic in laboratory animal experiments.

SELENIUM

Acute exposures can produce CNS effects, including nervousness, drowsiness, and convulsions, and eye and nasal irritation. Chronic exposure to selenium-containing compounds by inhalation can result in pallor, coated tongue, gastrointestinal disorders, nervousness, garlic breath, liver and spleen damage, anemia, and mucosal irritation. Discoloration, decayed teeth, skin eruptions, gastrointestinal distress, and loss of hair and nails have been reported in humans exposed orally. In livestock, excess intake can cause blind staggers (impaired vision, weak limbs, respiratory failure) and alkali disease (hair loss, sterility, atrophy of hooves, lameness, and anemia). Embryotoxic and teratogenic in animals. Selenium is carcinogenic in laboratory animals and may be anticarcinogenic and protective in humans. Selenium is an essential element. Its toxicity is related to chemical form.

TETRACHLOROETHENE (PERCHLOROETHYLENE)

Tetrachloroethene can depress the CNS and cause narcosis. It is irritating to mucous membranes and skin and can cause lung edema. Neurological effects on dry-cleaners have been reported. Chronic exposure may result in pathological changes in liver of laboratory animals. It may also affect the kidney. In humans, inhalation exposure may produce irritation of respiratory tract, nausea, headache, sleeplessness, and abdominal pains. Fatalities have been reported. It is carcinogenic in laboratory animals. An increased incidence of cancers among dry-cleaning workers exposed to several solvents has been described.

TOLUENE

Toluene is absorbed in humans following both inhalation and dermal exposure. At concentrations greater than 200 ppm (754 mg/ml), the primary acute effects are central nervous system depression and necrosis. At lower levels, nausea, fatigue, and incoordination have been reported. Chronic exposure to toluene vapors at concentrations of 200 to 800 ppm has been associated with central nervous and peripheral system effects, hepatomegaly, and hepatic and renal function changes.

TRICHLOROETHENE

Trichloroethene (TCE) is a central nervous system depressant after acute and chronic exposure. Oral exposures of humans to single doses of 15 to 25 ml (21 to 35 gm) of TCE have resulted in vomiting, abdominal pain, and transient unconsciousness. Absorption from the gastrointestinal tract is virtually complete and is proportional to concentrations and duration of exposure following inhalation. High level exposure can result in death from respiratory and cardiac failure. Industrial use and contact with TCE as a concentrated solvent is associated with adverse dermatological effects; however, no adverse effects are reported after exposure as a dilute, aqueous solution. Long-term inhalation exposure can affect liver and kidneys in animals. In humans, changes in liver enzyme have been associated with TCE exposure. TCE is classified as a Group B2 Carcinogen--Probable Human Carcinogen.

VINYL CHLORIDE

Acute occupational exposure to high concentrations of vinyl chloride can produce symptoms of narcosis. Respiratory tract irritation, bronchitis, headache, irritability, memory disturbances, and tingling sensations may also occur. Health effects associated with chronic occupational exposure include hepatitis-like liver changes, disturbances in visual and central nervous systems, decreased blood platelets and pulmonary function, and cardiovascular and gastrointestinal toxicity. Vinyl Chloride is classified as a Group A Known Human Carcinogen. Possible relationships between exposure and birth defects and fetal death have been reported. Chromosome aberrations have been reported in exposed workers.

XYLENES

The three xylene isomers, compounds having the same chemical constituents in a different configuration, have similar toxicological properties and are discussed together. Dermal absorption is reported to be minor after exposure to xylene vapor, but may be significant after contact with the liquid. Acute inhalation exposure in humans to high concentrations can depress the central nervous system and irritate mucous membranes. Changes in behavioral tests, manual coordination, balance, and electroencephalographic patterns have been reported in humans from chronic exposure; development of a tolerance to some of these effects has been reported.

ZINC

Zinc is an essential nutrient. The taste threshold to zinc is 15 ppm. Concentrations of soluble zinc salts of 40 ppm impart metallic taste. Acute adverse effects to inhalation

of zinc fumes include metal fume fever. Fever, nausea, vomiting, stomach cramps, and diarrhea may result from acute ingestions. Prolonged ingestion of zinc can result in irritability, muscular stiffness and pain, loss of appetite, and nausea. High levels of zinc in the diet may retard growth and produce defective mineralization of bone.

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NOTE: Health effects may be based on animal studies and do not imply that human exposure will have the same results.

Appendix J
INVESTIGATION-DERIVED WASTE MANAGEMENT

APPENDIX J Investigation-Derived Waste Management

TASK DESCRIPTION AND PURPOSE

During the SI, 29 drums were filled with generated waste that included drilling cuttings, sludges, groundwater, decontamination water, well development water, plastic, soiled personnel protective equipment, and other disposable items. Some drums contained a combination of the materials listed above. The investigation-derived waste sampling was implemented to reduce the quantity of waste requiring offsite disposal while also verifying that contaminated materials were disposed of properly.

According to "RCRA ARARs: Focus on Closure Requirements," in order for RCRA requirements to be applicable, a Superfund waste must be determined to be a listed or characteristic hazardous waste under RCRA. Sample results were evaluated based on characteristic level criteria only because no evidence exists that listed hazardous waste disposal occurred at the site.

Background samples were obtained and analyzed for the presence of organic and inorganic constituents. Samples were then obtained of borings from Sites 1 and 4 and analyzed for organic constituents. Previous use of the site did not indicate the need to analyze for the presence of inorganics. These samples showed no significant concentrations of organic constituents. Samples were obtained from Site 2 and analyzed for organics and inorganics. Six samples obtained from Site 2 were found to contain significant levels of Trichloroethene. Therefore, wastes from Site 2 will be disposed of as a hazardous waste.

At Site 1 (POL Facility), which contained six underground JP-4 storage tanks, samples were obtained and found to contain greater than 100 ppm total petroleum hydrocarbon (TPH) as deep as 8 feet. Soils containing greater than 100 ppm TPH must be remediated. Aeration of the contaminated soil is recommended.

The remaining drums of waste are suitable for disposal at a properly permitted solid waste landfill.

FIELD PROCEDURES

The disposal method selected for each investigation-derived waste material was based on field observations, field screening (HNu), and laboratory analyses of soil and water samples. All drums were sealed and labeled by boring and site and were stored in the Base hazardous waste generator storage area.

Appendix C contains the sample results pertaining to investigation-derived waste.

Appendix K
Geologist's Log
Well N-7
Alabama Air National Guard
Dannelly Field Municipal Airport

TABLE 4.—SAMPLE LOGS OF WELLS IN MONTGOMERY COUNTY, ALA,—Continued

TAB

YEA YEA

Thickness (feet)		Depth (feet)
Well N-2Continued		
angular, quartzose, and clay, reddish- een, and pale red-purple, ous	10	740
Clay, moderate reddish-brown, pale red-purple, and pale green, sandy, micaceous	10	750
ned, angular to subangular, quartzose, uginous; clay, varicolored, sandy; and greenish-gray, micaceous, pyritic	40	790
Well N-7		
(Samples described by John C. Scott) Owner: Alabama Air National Guard Driller: Acme D.	Drilling	S
le chalk:	· <u>·</u>	
Chalk, yellowish-gray, sity	121	333 133 133 133 133 133 133 133 133 133
gray, such, sugarcy measured and glauconitic gray, silty, slightly micaceous and glauconitic gray silty, slightly micaceous	14.	102
gray, silty, glauconitic, slightly mi-		
ly glauconitic and		195
sandy, slightly glauconitic and	2	77
	77	137
Sand, light gray, silty and fine- to medium-		
and micaceous, fossiliferous	22	159
ay, sirty and medium- to coarse-grained, subangular, quartzose, fossiliferous,		,
1	10	169
angular to subangular, quartzose, glauconitie	13	182
glay, sandy, dual cook, blancours, , micaceous alcarcours nt greenish gray, fine- to medium-grained,	14	196
lar to subangular, quartzose, very glau- ic, micaceous	9	202
gray, coarse-grained, angular to sub- quartzose, glauconitic	. 22	224
light gray, medium- to coarse-grained, ilar to subangular, quartzese, glauconitic,	ę	į
and sandstone, white, carcalcons Sand, light gray, coarse-grained, angular to sub-	3	1 67
ular, quartzose, sparsay glaurovitie vollowish aray medium, grained ang		258
ngular, quartzose, sparsely glauconitic light gray, medium- to coarse-grained,	ਜ਼ ਂ	279
angular to subangular, quartzose, sparsely glauconitie, and micaceous	12	291

GY, ND- R RESCRICES, mon'TGOMENT COUNTY

TABLE 4.—SAMPLE LOGS OF WELLS IN MONTGOMERY COUNTY, ALA.—Continued

Thickness (feet)	Depth (feet)	l
Well N-7-Continued		1
Sand, light greenish-gray, medium-grained, angular to subangular, quartzose, glauconitic, micaceous, and clay, gray, sandy, micaceous	301	
glauconitic, sparsely micaceous Sand, light greenish-gray, silty and fine- to medium-grained, angular to subangular	313	
quartzose, glauconitic, micaceous	344	
quartzose, sparsely glauconitic, and micaceous 46 Sand, light gray, medium- to coarse-grained, angular to subangular, quartzose, sparsely glauconitic	390	
and micaceous 37 Sand, light gray, coarse-grained, angular to subangular	427	
quartzose, sparsely glauconitic and micaceous Sand, light gray, medium- to coarse-grained, angular to subangular, quartzose, sparsely glauconitic mica-	447	
ceous and pyritic, and clay, gray, sandy, micaceous 10 Sand, light gray, coarse-grained, angular to subangular.	457	
quartzose, sparsely glauconitic and micaceous13	470	
(Samples described by H. L. Reade, Jr.) Owner: U. S. Geological Survey (test well GS-9) Driller: Black Relt Drilling Co		
	20 160	
Sand, light greenish-gray, coarse-grained, sub- angular to subrounded, frosted, quartzose, glau- conitic, and sandstone, light greenish-gray, cal-		
careous, glauconitic	170	
Sand, light greenish-gray, medium-grained, sub- angular to subrounded, frosted, quartzose, glauconitic, fossiliferous, and clay, greenish-	180	
gray, sandy, glauconitic, fossiliferous	220	
rounded, slightly frosted, quartzose, glauconitic 30 Sand, light greenish-gray, medium- to coarse-grained.	230	
subangular to subrounded, quartzose, glauconitic Sand, light greenish-gray, medium- to coarse-grained, subangular to subrounded, slightly frosted, quartzose.	240	
glauconitic, and clay, greenish-gray, micaceous Sand, light greenish-gray, coarse-grained, subangular to subrounded, frosted, quartzose, glauconitic,	270	
fossiliterous 40 Clay, greenish-gray, micaceous, and sand, light greenish-	310	